GROUNDWATER MONITORING REPORT SECOND QUARTER 2006 PACIFIC AIRMOTIVE CORPORATION 2940 AND 3003 NORTH HOLLYWOOD WAY BURBANK, CALIFORNIA



Prepared for:

LOCKHEED MARTIN







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August 21, 2006

Ms. Rachel Loftin

Remedial Project Manager

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Dear Rachel:

Enclosed please find one (1) copy of the Groundwater Monitoring Report, Second Quarter 2006, Pacific Airmotive Corporation, 2940 and 3003 North Hollywood Way, Burbank, California. Please do not hesitate to contact me if you have any questions or comments.

Regards,

Lisa A. Hamilton

Manager, MidAtlantic/Southeast/Western Regions

cc Linda Gertler, LMC (w/out enclosure)
Ken Martins, CH2M Hill (with enclosure)
Dixon Oriola, LARWQCB (with enclosure)
Alex Lapostol, E2 (with enclosure)

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August 2006

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SECTION I INTRODUCTION

1.0 INTRODUCTION

On behalf of Lockheed Martin Corporation (LMC), Tetra Tech, Inc. (Tetra Tech) has prepared this groundwater monitoring report for two Pacific Airmotive Corporation (PAC) properties within the Burbank Operable Unit (BOU) in Burbank, California (Figure 1-1). LMC is performing work requested by the U.S. Environmental Protection Agency (EPA) in a letter directed to General Electric (GE) dated October 20, 2005 due to a settlement agreement between PAC, an indirect wholly-owned subsidiary of GE, and LMC.

In the October 20, 2005 letter, the EPA requested GE to initiate four quarters of groundwater sampling of the eight (8) existing wells at the PAC properties based on previous facility operations, detection of constituents, lack of current groundwater results, and recent regulatory concerns related to potential sources associated with emergent chemicals within the BOU. The EPA required analysis of the groundwater for volatile organic compounds (VOCs), 1,2,3-trichloropropane (1,2,3-TCP), Title 22 metals, including thallium and dissolved (total) chromium, hexavalent chromium, 1,4-dioxane, N-Nitrosodimethylamine (NDMA), perchlorate, nitrate/nitrite, common cations and anions, dissolved oxygen, sulfide, and dissolved iron and manganese.

1.1 SITE LOCATION AND DESCRIPTION OBJECTIVE

The PAC properties are located at 2940 and 3003 North Hollywood Way within the north-central portion of the BOU (Figure 1-1). The property at 2940 North Hollywood Way was identified as the Main Facility, and the property at 3003 North Hollywood Way was identified as the Jet Engine Test Cell Facility. Both facilities were historically associated with the manufacturing, design, and repair of aircraft and aircraft engines. Structures located on both PAC properties are currently vacant.

1.2 OBJECTIVE

The purpose of this groundwater monitoring report is to comply with the provisions of the EPA October 20, 2005 letter. The objective of this monitoring report is to present

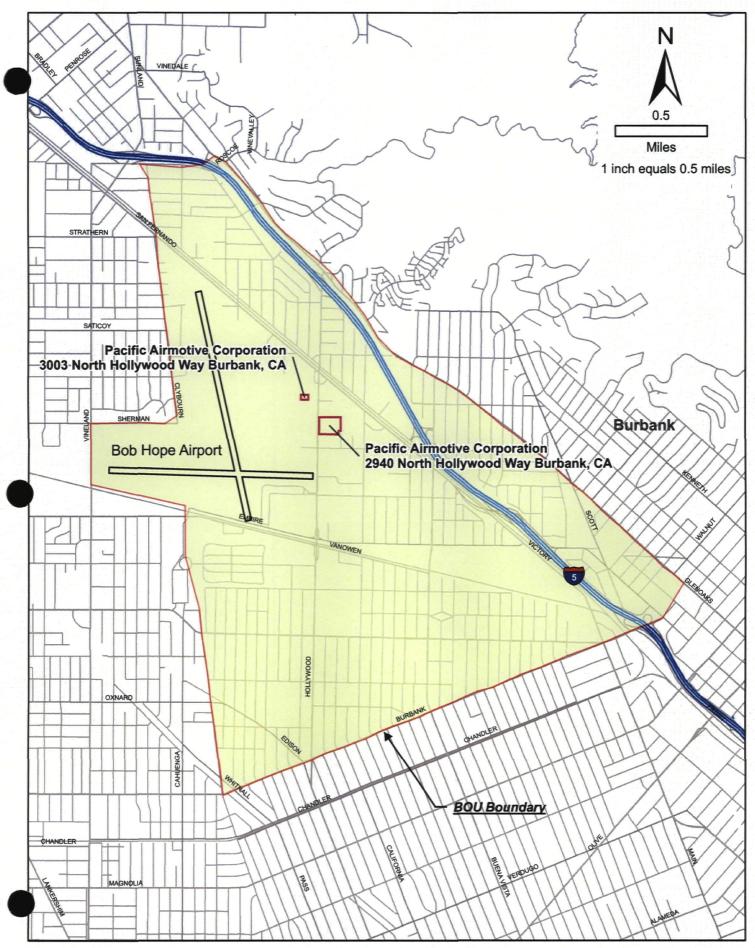


Figure 1-1 - BOU Boundary Map

groundwater data collected during the second quarter 2006. The groundwater data is being collected to assist the EPA in assessing the current groundwater quality and conditions at the above mentioned monitoring wells and within the BOU. The quarterly monitoring report presents field, laboratory analytical results, and quality control data collected during groundwater level and water quality monitoring.

1.3 REPORT ORGANIZATION

The second quarter 2006 quarterly groundwater monitoring report has been organized into the following six (6) sections:

- Section 1. <u>Introduction</u>: introduces the project and presents the objectives and report format.
- Section 2. <u>Subsurface Conditions</u>: presents the site geologic and hydrogeologic setting.
- Section 3. <u>Description of Historical Areas of Concern</u>: identifies the areas of groundwater concern beneath the PAC properties.
- Section 4. <u>Groundwater Monitoring Procedures</u>: summarizes the groundwater monitoring activities, groundwater measurements, and laboratory analysis conducted.
- Section 5. <u>Groundwater Analytical Results</u>: discusses groundwater monitoring results.
- Section 6. <u>References</u>: lists the references used to prepare this quarterly groundwater monitoring report.

SECTION 2 SUBSURFACE CONDITIONS

2.0 SUBSURFACE CONDITIONS

2.1 GEOLOGY

The PAC properties are located in the southeastern portion of the San Fernando Valley (SFV) between the Santa Monica and Verdugo mountains. The SFV is located on the northwestern block of the Los Angeles Basin within the Transverse Ranges Geomorphic Province, an east-west trending unit composed of subparallel ranges separated by alluviated, synclinal valleys and prominent faults. The SFV is bordered to the north by the Santa Susana and San Gabriel mountains, to the east by the Verdugo Mountains, to the south by the Santa Monica Mountains, and to the west by the Simi Hills. These uplands are composed of crystalline bedrock of Precambrian to Mesozoic in age and sedimentary units from Cretaceous to Pleistocene in age. The crystalline bedrock and sedimentary units were eroded from the uplands during the Quaternary Period and deposited as more than 2,000 feet of alluvium in the SFV. The only major structural feature within close proximity to the PAC properties is the Verdugo Fault, which is approximately one mile to the northeast and trends northwesterly along the base of the Verdugo Mountains (Tetra Tech, 2006a).

2.2 HYDROGEOLOGY

The PAC properties are located within the San Fernando Basin (SFB), one (1) of four (4) distinct groundwater basins that encompass the entire watershed of the Los Angeles River and its tributaries within the SFV (also known as the Upper Los Angeles River Area – ULARA). Groundwater within the eastern portion of the SFB flows mainly through two sedimentary units: 1) Older Alluvium of Pleistocene age and 2) Younger Alluvium of Holocene age. The Older Alluvium is comprised of sand, gravel, and boulders in the northwestern portion of the BOU to interbedded silt and sand in the eastern and southern portions of the BOU. The Younger Alluvium is comprised of coarse sand, gravel, and cobbles interbedded with finer-grained units of sand, silty sand, sandy silt, silty clay, and minor gravelly sand. Groundwater flow within the Older Alluvium has been observed to be locally semi-confined to confined. The Younger alluvium is generally unconfined to

semi-confined, depending upon the location and thickness of fine grained interbeds (Tetra Tech, 2006a).

The aquifer in the Younger Alluvium within the BOU has been divided into five hydrostratigraphic units (HSU) based on electrical resistivity responses in geophysical logs (Hargis & Associates, 1991; Simon Hydro-Search, 1993). The five HSUs of the Younger Alluvium are identified from upper to lower as A', X, A, Y, and B. The A', A, and B units are generally composed of coarser-grained material (coarse sands, gravels, and cobbles). The X and Y HSUs separate the three (A', A, B) HSUs listed above and consist of relatively finer-grained material including sand, silty sand, and silt. Based on the stratigraphic position of the units and the groundwater gradient, the A', X, or A HSU may locally represent water table conditions depending on geographic location within the project area.

Groundwater flow direction in the SFB is generally toward the southeast. Groundwater velocities in the BOU range from approximately 300 to 900 feet per year (ULARA, 2005).

SECTION 3 DESCRIPTION OF HISTORICAL AREA OF CONCERN

3.0 DESCRIPTION OF HISTORICAL AREAS OF CONCERN

After reporting a jet fuel spill to the Los Angeles Regional Water Quality Control Board (LA-RWQCB) in 1987, PAC agreed to install MW-1 and MW-2 at the Jet Engine Test Facility downgradient of the location of the fuel spill. In 1992, in an effort by the LA-RWQCB to assess the groundwater analytes underlying the PAC properties, monitoring well MW-3 was installed at the Jet Engine Test Cell Facility, and wells MW-4 through MW-8 were installed at the Main Facility (Figure 3-1).

The EPA issued a Unilateral Administrative Order (UAO) in 1994 which required PAC to perform soil and groundwater investigations. As part of the soil investigation, PAC conducted soil gas surveys across the PAC properties to assess the nature and extent of vapor and non-vapor phase analytes in the unsaturated zone. Since 1997, when PAC became an indirect wholly owned subsidiary of GE, PAC, through GE technical and legal representatives acting on its behalf, has been working with the LA-RWQCB to further investigate and remediate PAC properties (Tetra Tech, 2006a).

Semi-annual groundwater monitoring from June 1987 through December 1988 indicated elevated levels of trichloroethene (TCE) and tetrachloroethene (PCE) in monitoring wells MW-1 and MW-2 (Table 3-1). Groundwater monitoring from September 1992 through January 1995 showed PCE and TCE concentrations exceeding regulatory maximum contaminant levels (MCLs) of 5 micrograms per liter (μ g/L) in wells MW-3 through MW-8 (Table 3-2). Monitoring wells MW-1 and MW-2 were both dry during this time period.



Figure 3-1 - PAC Wells Location Map

Table 3-1 Historical Analysis From 1987 – 1989 (Reported in μ g/L)

	6/1	8/87	12/2	9/87	6/1	4/88	12/15/88		
	PCE	TCE	PCE	TCE	PCE	TCE	PCE	TCE	
MCL	5	5	5	5	5	5	5	5	
Composite of MW-1 & MW-2	130	32							
MW-1	130*	32*	67	24	160	31	75	12	
MW-2	130*	32*	190	41	200	33	130	15	

Notes:

All concentrations in µg/L

**Bold - Result above MCL

**Result based on composite sample

Table 3-2 Historical Analysis From 1992 - 1995 (Reported in µg/L)

Well ID	9/15-	16/92	12/16-19/92		7/19-	20/94	12/25	-26/94	1/30-31/95	
Well ID	PCE	TCE	PCE	TCE	PCE	TCE	PCE	TCE	PCE	TCE
MCL	5	5	5	5	5	5	5	5	5	5
MW-1	dry	dry	dry	dry	dry	dry	dry	dry	dry	dry
MW-2	dry	dry	dry	dry	dry	dry	dry	dry	dry	dry
MW-3	39	11	47	12	18	6.4	58	8.8	63	7.8
MW-4	460	46	400	41	22	6.3	25	3.6	13	2.2
MW-5	2100	440	64	13	40	8.9	150	24	49	6.9
MW-6	910	250	490	120	39	7.4	1300	170	800	110
MW-7	87	18	420	49	43	11	2000	88	490	19
MW-8	1700	160	1200	94	21	5.1	1800	170	1800	130

Notes:

All concentrations in μg/L **Bold** – Result above MCL

SECTION 4
GW MONITORING
PROCEDURES

4.0 GROUNDWATER MONITORING PROCEDURES

4.1 GROUNDWATER LEVEL MEASUREMENTS

Water levels in the eight (8) monitoring wells were measured on June 2 and June 6, 2006 using a water level meter consisting of a liquid sensor attached to a measuring tape that was lowered down into the well until water was encountered. Water level measurements were recorded on well purging forms (Appendix A) and are presented in Table 4-1. Groundwater monitoring wells MW-1 and MW-2 were dry. Groundwater elevation contours are shown on Figure 4-1. The groundwater flow direction is to the east.

Table 4-1
Summary of Groundwater Elevations

Well Number	HSU	Date Measured	Top of Casing (TOC) Elevation (feet msl)	Groundwater Depth from TOC (feet)	Groundwater Elevation (feet msl)
MW-1	NA	dry	NA	dry	dry
MW-2	NA	dry	NA	dry	dry
MW-3	NA	06/02/2006	NA*	241.56	NA
MW-4	A	06/06/2006	700.5	227.61	472.89
MW-5	A	06/06/2006	704.2	229.67	474.53
MW-6	A	06/06/2006	701.2	227.96	473.24
MW-7	A	06/06/2006	696.4	225.61	470.79
MW-8	A	06/06/2006	NA*	227.25	NA

Note:

HSU - Hydrostatic unit

TOC - Top of casing

msl - Mean sea level

NA - Not available

^{*} TOC will be surveyed prior to the third quarter 2006 monitoring



Figure 4-1 - Second Quarter 2006 WT HSU's Groundwater Elevation, PAC

4.2 WELL PURGING AND SAMPLING

Well development was completed in March 2006 in order to optimize groundwater production within each well prior to the initial quarterly sampling. This was done because groundwater monitoring and sampling had not been completed at the eight (8) PAC wells since 1995.

Prior to collecting the groundwater samples in June 2006, a minimum of three well volumes was purged from monitoring wells MW-3 through MW-8 using a submersible pump. Water temperature, pH, conductivity, dissolved oxygen, and turbidity were measured throughout the purging process using a field water quality monitoring system. Stabilization of these parameters served as an indication of water representative of the formation, and their values were recorded on well purging forms (Appendix A).

The groundwater samples were collected using a down-hole submersible pump for monitoring wells MW-3 through MW-8. Groundwater samples were collected from a nozzle attached to the pump hose and placed directly into sample containers provided by the laboratory. Decontamination procedures were followed after each monitoring well was sampled to avoid cross-contamination between wells. The water samples were placed on ice in a cooler to maintain a temperature of ⁺/. 4°C pending delivery to Calscience Environmental Laboratories, Inc., a State of California certified laboratory, for analysis. A completed chain-of-custody form accompanied the shipment of samples to the laboratory to ensure accountability for the samples from the time of collection to the time of analysis.

4.3 LABORATORY ANALYSIS

Groundwater samples were collected from the six (6) groundwater monitoring wells (MW-3 through MW-8) on June 19 and 20, 2006 at the PAC facility. Samples analyzed for dissolved metals were filtered in the field using a disposable filter.

The EPA has requested that groundwater samples from the PAC wells be analyzed for specific constituents using analytical methods consistent with those of the BOU groundwater sampling events as follows:

- ➤ VOCs, including MTBE, using EPA Method 8260B;
- > 1,2,3-TCP, using EPA Method 504.1 or 524M;
- > Title 22 metals, including thallium and dissolved (total) chromium, using EPA Method 6010B/7470A;
- > Hexavalent chromium, using EPA Method 218.6;
- > 1,4-dioxane, using EPA Method 8270C (M);
- > NDMA, using EPA Method 1625C (M);
- > Perchlorate, using EPA Method 314.0;
- ➤ Nitrate/nitrite, using EPA Method 300.0;
- > Cations, using EPA Method 6010B;
- > Anions, using EPA Method 300.0;
- Dissolved oxygen, using EPA Method SM 4500-O G;
- > Sulfide, using EPA Method 376.2; and
- Dissolved iron and manganese using EPA 3005A Filter/EPA 200.8.

SECTION 5 GW ANALYTICAL RESULTS

5.0 GROUNDWATER ANALYTICAL RESULTS

Based on the data collected during the second quarter 2006 groundwater sampling event, compounds are reported and compared to their respective MCL or California drinking water notification level (CDWNL). The MCL or CDWNL concentrations are based on the lowest value in "A Compilation of Water Quality Goals, California Regional Water Quality Control Board, Central Valley Region," dated September 2004. Copies of the laboratory analytical data reports are included in Appendix B. A summary of the analytical detected is provided in Tables 5-1 through 5-6. A summary of the analytical results is presented in the following subsections.

5.1 VOC ANALYTICAL RESULTS

Groundwater samples collected from six (6) groundwater monitoring wells were analyzed for VOCs. A summary of the analytical results are presented in Table 5-1 and discussed below:

- Acetone was detected in one (1) groundwater sample (MW-8) at a concentration of 10 μg/L and was estimated in three (3) groundwater samples (MW-3, MW-5, and MW-7) with concentrations of 9.7 μg/L, 6.3 μg/L, and 9.8 μg/L, respectively. However, these three estimated sample results are considered not to have originated from the native sample. Laboratory cross-contamination (shown in the method blank) likely caused these detections.
- > Carbon Tetrachloride was detected in five (5) groundwater samples (MW-3, MW-4, MW-5, MW-6, and MW-8) with concentrations of 1.1 μg/L, 0.81 μg/L, 2.4 μg/L, 3.1 μg/L, and 0.79 μg/L, respectively.
- > Chloroform was detected in five (5) groundwater samples (MW-3, MW-4, MW-5, MW-6, and MW-8) with concentrations of 1.4 μg/L, 1.3 μg/L, 2.0 μg/L, 2.4 μg/L, and 1.1 μg/L, respectively, and was estimated in one (1) groundwater sample (MW-7) with a concentration of 0.71 μg/L.
- > Methylene Chloride was estimated in all six (6) groundwater samples (MW-3 through MW-8) with concentrations of 3.6 μg/L, 4.5 μg/L, 4.6 μg/L, 3.9 μg/L, 3.6 μg/L, and 4.2 μg/L, respectively. However, these detections are considered not to have originated from the native sample. Laboratory cross contamination (shown in the method blank) likely caused the detections.
- > Tetrachloroethene was detected in all six (6) groundwater samples (MW-3 through MW-8) with concentrations of 36 μg/L, 120 μg/L, 150 μg/L, 150 μg/L, 50 μg/L, and 160 μg/L, respectively.

- > 1,2-Dichloroethane was detected in one (1) groundwater sample (MW-8) at a concentration of 1.6 μg/L and was estimated in two (2) groundwater samples (MW-3 and MW-5) with concentrations of 0.46 μg/L and 0.29 μg/L, respectively.
- > 1,1-Dichloroethene was detected in three (3) groundwater samples (MW-3, MW-5, and MW-6) with concentrations of 2.6 μg/L, 2.6 μg/L, and 3.4 μg/L, respectively, and estimated in three (3) groundwater samples (MW-4, MW-7, and MW-8) with concentrations of 0.64 μg/L, 0.41 μg/L, and 0.97 μg/L, respectively.
- > 1,1,2-Trichloro-1,2,2-trifluoroethane was estimated in five (5) groundwater samples (MW-3, MW-4, MW-5, MW-6, and MW-8) with concentrations of 1.9 μg/L, 1.1 μg/L, 1.6 μg/L, 1.1 μg/L, and 2.2 μg/L, respectively.
- Arr Trichloroethene was detected in all six (6) groundwater samples (MW-3 through MW-8) with concentrations of 12 μg/L, 48 μg/L, 84 μg/L, 75 μg/L, 17 μg/L, and 58 μg/L, respectively.

A review of the VOC analytical data shows that four (4) compounds were detected above their respective MCL. Carbon tetrachloride was detected above the MCL of $0.5~\mu g/L$ in five (5) groundwater samples (MW-3 through MW-6 and MW-8). Tetrachloroethene was detected above the MCL of $5~\mu g/L$ in all six (6) groundwater samples. The compound 1,2-dichloroethane was detected above the MCL of $0.5~\mu g/L$ in one (1) groundwater sample (MW-8). Trichloroethene was detected above the MCL of $5~\mu g/L$ in all six (6) groundwater samples.

Table 5-1 Summary of Detected VOCs Analytical Results EPA Method 8260B (Reported in µg/L)

Well ID	Acetone	Bromo-dichloromethane	Carbon Tetrachloride	Chloroform	Chloromethane	Methylene Chloride	Tetrachloroethene	1,2-Dichloroethane	1,1-Dichloroethene	1,1,2-Trichloro-1,2,2- trifluoroethane	Trichloroethene
MCL	NA	80 ¹	0.5	80 ¹	NA	NA	5 ²	0.5 ²	6 ²	1,200 ²	5 ²
MW-3	9.7 ^{J,B}	<0.27	1.1	1.4	<1.8	3.6 ^{J,B}	36	0.46 ^J	2.6	1.9 ^J	12
MW-4	<6.1	<0.27	0.81	1.3	<1.8	4.5 ^{J,B}	120	<0.22	0.64 ^J	1.0 ^J	48
MW-4D	9.5 ^J	< 0.27	0.88	1.2	<1.8	2.9 ^{J,B}	120	<0.22	0.82 ^J	1.1 ^J	48
MW-5	6.3 ^{J,B}	0.29 ^J	2.4	2.0	<1.8	4.6 ^{J,B}	150	0.29 ^J	2.6	1.6 ^J	84
MW-6	<6.1	<0.27	3.1	2.4	<1.8	3.9 ^{J,B}	150	<0.22	3.4	1.1	75_
	10	-0.27	<0.42	0.71	<1.8	3.6 ^{J,B}	50	<0.22	0.41 ^J	<0.54	17
MW-7	9.8 ^{J,B}	<0.27	10.42	0.71		4.2 ^{J,B}					

Note:

- 1 US EPA MCL
- ² California Primary MCL
- Analyte was present in the associated method blank and the result is considered not to have originated from the native sample.
- Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated

MCL = Maximum contaminant level

NA = Not available

5.2 EMERGENT CHEMICALS ANALYTICAL RESULTS

Groundwater samples collected from the six (6) groundwater monitoring wells were analyzed for 1,4-dioxane, NDMA, and 1,2,3-TCP. A summary of the analytical results is presented in Table 5-2 and discussed below:

- \triangleright 1,4-Dioxane was not detected in any of the groundwater samples above the laboratory method detection limit of 0.40 μ g/L.
- > NDMA was not detected in any of the groundwater samples above the laboratory method detection limit of $0.00048 \mu g/L$.
- > 1,2,3-TCP was detected in all six (6) groundwater samples with concentrations ranging from 0.016 ug/L (MW-7) to 0.19 μg/L (MW-6).

A review of the emergent chemical analytical data reveals that 1,2,3-TCP was detected above its CDWNL of $0.005 \mu g/L$ in all six (6) groundwater samples.

Table 5-2
Emergent Chemicals Analytical Results
(Reported in µg/L)

Well ID	1,4-Dioxane by EPA Method 8270 SIM	NDMA by EPA Method 1625C(M)	1,2,3-TCP by EPA Method 504.1
CDWNL	3	0.01	0.005
MW-3	<0.40	<0.00048	0.044
MW-4	< 0.40	<0.00048	0.028
MW-4D	<0.40	<0.00048	0.027
MW-5	<0.40	<0.00048	0.14
MW-6	<0.40	<0.00048	0.19 ^B
MW-7	<0.40	<0.00048	0.016
MW-8	<0.40	<0.00048	0.068

Note:

CDWNL = California Drinking Water Notification Level

NDMA = N-Nitrosodimethylamine 1,2,3-TCP = 1,2,3-Tricloropropane

MCL = Maximum contaminant level

B Analyte was present in the associated method blank.

5.3 DISSOLVED IRON AND MANGANESE ANALYTICAL RESULTS

Groundwater samples collected from six (6) groundwater monitoring wells were analyzed for dissolved iron and manganese. A summary of the analytical results are presented in Table 5-3 and discussed below.

- ➤ Dissolved Iron was estimated in all six (6) groundwater samples with concentrations ranging from 0.0172 mg/L (MW-3) to 0.0342 mg/L (MW-4).
- ➤ **Dissolved Manganese** was detected in all six (6) groundwater samples with concentrations ranging from 0.00184 mg/L (MW-7)to 0.00497 mg/L (MW-4).

A review of the dissolved iron and manganese analytical results reveal that groundwater samples did not contain concentrations that exceeded their respective MCL.

Table 5-3
Dissolved Metals Analytical Results
EPA Method 6010B/7470A
(Reported in mg/L)

Well ID	Iron	Manganese
MCL	0.3 1	0.05 1
MW-3	0.0172 ^J	0.00432
MW-4	0.0342 ^J	0.00497
MW-4D	0.0246 ^J	0.00488
MW-5	0.0301 ^J	0.00261
MW-6	0.0260 ^J	0.00439
MW-7	0.0210 ^J	0.00184
MW-8	0.0196 ^J	0.00235

Note:

- 1 US EPA MCL
- ^B Analyte was present in the associated method blank.
- Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated.

MCL = Maximum Contaminant Level.

5.4 INORGANIC ANALYTICAL RESULTS

Groundwater samples collected from six (6) groundwater monitoring wells were analyzed for hexavalent chromium, perchlorate, anions, dissolved oxygen and other constituents. A summary of the analytical results are presented in Table 5-4 and discussed below:

- ➤ Hexavalent Chromium was detected in all six (6) groundwater samples at concentrations ranging from 0.0014 mg/L (MW-7) to 0.0029 mg/L (MW-6).
- ➤ Chloride was detected in all six (6) groundwater samples with concentrations ranging from 37 mg/L (MW-5) to 41 mg/L (MW-3).
- > Nitrite was not detected in any of the groundwater samples above the laboratory reporting limit.
- ➤ Nitrate was detected in all six (6) groundwater samples with concentrations ranging from 10 mg/L (MW-7) to 16 mg/L (MW-4).
- > Sulfate was detected in all six (6) groundwater samples with concentrations ranging from 67 mg/L (MW-7) to 75 mg/L (MW-5).
- > Sulfide was not detected in any of the groundwater samples above the laboratory reporting limit.
- ➤ **Perchlorate** was estimated in all six (6) groundwater samples with concentrations ranging from 0.00056 mg/L (MW-6) to 0.00099 mg/L (MW-5).

➤ **Dissolved Oxygen** was detected in all six (6) groundwater samples with concentrations ranging from 7.41 mg/L (MW-3) to 8.17 mg/L (MW-6).

A review of the inorganic analytical data reveals that nitrate equaled or exceeded the California drinking water notification level of 10 mg/L. Hexavalent chromium was not detected above the regulatory action level of 0.05 mg/L in all six (6) groundwater samples.

Table 5-4
Inorganic Analytical Results
(Reported in mg/L)

Well ID	Hexavalent Chromium	Chloride	Nitrite	Nitrate	Sulfate	Sulfide	Perchlorate	Dissolved Oxygen
Regulatory Action Level	0.05	250 ²	1 ²	10 ²	250 ²	NA	0.006 ³	NA
MW-3	0.0016 ^B	41	<0.015	13	71	<0.042	0.00075 ^J	7.41
MW-4	0.0015 ^B	40	<0.015	16	72	<0.042	0.00058 ^J	7.89
MW-4D	0.0015 ^B	44	<0.015	12	73	<0.042	0.00071 ^J	7.82
MW-5	0.0017 ^B	37	<0.015	13	75	< 0.042	0.00099 ^J	7.36
MW-6	0.0029 ^B	38	< 0.015	13	74	< 0.042	0.00056 ^J	8.17
MW-7	0.0014 ^B	38	< 0.015	10	67	< 0.042	0.00059 ^J	7.71
MW-8	0.0015 ^B	40	<0.015	12	73	< 0.042	0.00076 ³	7.54

Note:

- 1 Hexavalent chromium currently regulated using MCL for total chromium
- ² California Secondary MCL
 - ³ California Drinking Water Notification Level
 - Analyte was present in the associated method blank. However, the method blank concentrations are below the EPA validation criteria for hexavalent chromium. Report value is not affected.
 - J Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated.

NA = Not available.

5.5 CATION ANALYTICAL RESULTS

Groundwater samples collected from six (6) groundwater monitoring wells were analyzed for cations. A summary of the analytical results is presented in Table 5-5 and discussed below:

➤ Calcium was detected in all six (6) groundwater samples (MW-3 through MW-8) at concentrations ranging from 98.3 mg/L (MW-8) to 112 mg/L (MW-6).

- ➤ Magnesium was detected in all six (6) groundwater samples (MW-3 through MW-8) with concentrations ranging from 29.9 mg/L (MW-5) to 34.4 mg/L (MW-7).
- ➤ Potassium was detected in all six (6) groundwater samples (MW-3 through MW-8) with concentrations ranging from 5.45 mg/L (MW-7) to 5.87 mg/L (MW-3 and MW-6).
- > Sodium was detected in all six (6) groundwater samples (MW-3 through MW-8) with concentrations ranging from 36.6 mg/L (MW-5) to 38.8 mg/L (MW-3).

California drinking water notification levels and MCLs have not been established for cations.

Table 5-5
Cations Analytical Results
EPA Method 6010
(Reported in mg/L)

Well ID	Calcium	Magnesium	Potassium	Sodium
MW-3	107	32.7	5.87 ^B	38.8 ^B
MW-4	104 ^E	33.7	5.56	37.9
MW-4D	104	33.4 ^B	5.41	37.4
MW-5	109	29.9 ^B	5.66	36.6
MW-6	112	30.8 ^B	5.87	37.0
MW-7	98.3	34.4	5.45 ^B	37.9 ^B
MW-8	104	33.3	5.74 ^B	37.7 ^B

Note:

^B Analyte was present in the associated method blank

^E Concentration exceeds the calibration range

5.6 TITLE 22 METAL ANALYTICAL RESULTS

Groundwater samples collected from six (6) groundwater monitoring wells were analyzed for Title 22 metals. A summary of the analytical results are presented in Table 5-6 and only metal analytes detected above the laboratory reporting limit are listed below:

- **Barium** was detected in all six (6) groundwater samples with concentrations ranging from 135 μ g/L (MW-5) to 146 μ g/L (MW-4 and MW-7).
- > Cadmium was estimated in one (1) groundwater sample (MW-6) with a concentration of 1.22 μg/L.

- > Chromium was detected in all six (6) groundwater samples (MW-3 through MW-8) with concentrations ranging from 5.61 μg/L (MW-4) to 7.93 μg/L (MW-6).
- > Cobalt was estimated in three (3) groundwater samples (MW-3, MW-7, and MW-8) with concentrations of 1.97 μg/L, 1.71 μg/L and 1.94 μg/L, respectively.
- Nickel was estimated in three (3) groundwater samples (MW-3, MW-7, and MW-8) with concentrations of 3.65 μ g/L, 3.38 μ g/L and 2.39 μ g/L, respectively.
- > Selenium was estimated in three (3) groundwater samples (MW-4D, MW-5, and MW-6) with concentrations of 8.31 μg/L, 3.93 μg/L and 3.75 μg/L, respectively.
- > Thallium was estimated in four (4) groundwater samples (MW-3 through MW-6) with concentrations ranging from 3.21 μg/L (MW-3) to 11.2 μg/L (MW-6).
- > Vanadium was estimated in all six (6) groundwater samples with concentrations ranging from 3.26 μg/L (MW-3) to 4.77 μg/L (MW-4).
- > Zinc was detected in all six (6) groundwater samples with concentrations ranging from 85.0 μg/L (MW-4) to 181 μg/L (MW-6).

A review of analytical results for metals shows that they are all below their respective MCL with exception of thallium.

Table 5-6 **Title 22 Metals Analytical Results** EPA Method 6010B/7470A (Reported in $\mu g/L$)

Well ID	Antimony	Arsenic	Barium	Beryllium	Cadmium	Total Chromium	Cobalt	Copper	Lead	Mercury	Molybdenum	Nickel	Selenium	Silver	Thallium	Vanadium	Zinc
MCL	6	10	1000	4	5	50	NA	1000	15	2	NA	100	50	100	2	NA	5,000
MW-3	<2.09	<3.08	141	<0.176	<0.350	7.41 ^B	1.97 ^{J,B}	<1.34	<2.36	<0.0672	<0.800	3.65 ^{J,B}	<2.95	<0.400	3.21	3.26 ^J	110
MW-4	<2.09	<3.08	146	<0.176	<0.350	5.61	<0.696	<1.34	<2.36	<0.0672	<0.800	<1.37	<2.95	<0.400	6.94 ³	4.773	85.0
MW-4D	<2.09	<3.08	145	< 0.176	<0.350	6.22	<0.696	<1.34	<2.36	<0.0672	<0.800	<1.37	8.31 ^{J,B}	<0.400	5.94 ^{J,B}	4.67 ³	122
MW-5	<2.09	<3.08	135	<0.176	<0.350	6.48	<0.696	<1.34	<2.36	<0.0672	<0.800	<1.37	3.93 ^{J,B}	<0.400	4.98 ^{J,B}	3.53 ^J	144
MW-6	<2.09	<3.08	143	<0.176	1.22	7.93	<0.696	<1.34	<2.36	<0.0672	<0.800	<1.37	3.75 ^{J,B}	<0.400	11.2 ^{J,B}	4.31 ^J	181
MW-7	<2.09	<3.08	146	<0.176	<0.350	6.41 ^B	1.71 ^{J,B}	<1.34	<2.36	<0.0672	<0.800	3.38 ^{J,B}	<2.95	<0.400	<2.33	3.93 ^J	110
MW-8	<2.09	<3.08	145	<0.176	<0.350	6.50 ^B	1.94 ^{J,B}	<1.34	<2.36	<0.0672	<0.800	2.39 ^{J,B}	<2.95	<0.400	<2.33	3.50 ^J	93.7

Note:

MCL = Maximum contaminant level

NA = Not available

^{*} California Drinking Water Notification Level

Analyte was present in the associated method blank

Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated.

5.7 DATA VERIFICATION AND VALIDATION

Laboratory data underwent verification and validation including laboratory control samples (LCS), matrix spike duplicates (MSD), and method blanks. All samples received by the laboratory were analyzed within holding times specified by USEPA SW-846. Appendix C presents a summary of the quality control and quality assurance (QA/QC).

SECTION 6 REFERENCES

6.0 REFERENCES

- California Environmental Protection Agency Regional Water Quality Control Board Central Valley Region, 2004. A Compilation of Water Quality Goals. September 2004.
- Hargis & Associates, 1991. Installation of Groundwater Monitor Wells Along Vanowen Street, Lockheed Engineering and Science Company, Burbank, California.
- Simon Hydro-Search, 1993. Phase I Final Remedial Design Report, Burbank Operable Unit, Vols. V & VI prepared for LESAT, September 30, 1993.
- Tetra Tech, 2006a. Groundwater Monitoring and Sampling Work Plan, Pacific Airmotive Corporation Properties Located at 2940 and 3003 North Hollywood Way, Burbank, California, March 2006.
- Tetra Tech, 2006b. Groundwater Monitoring Report, First Quarter, Pacific Airmotive Corporation, 2940 and 3003 North Hollywood Way, Burbank, California, May 2006.
- U.S. Environmental Protection Agency, October 20, 2005. Letter to General Electric Company.
- Watermaster (ULARA), 2005. Watermaster Service in the Upper Los Angeles River Area Los Angeles County 2003 2004 Water Year, Upper Los Angeles River Area, May 2005.

APPENDIX A FIELD DATA LOG SHEETS

APPENDIX A

FIELD DATA LOG SHEETS

TABLE 1 List of Groundwater Monitoring Wells for for the 2006 Groundwater Monitoring Program
Burbank Operable Unit
WATER LEVELS MEASUREMENT FOR PAC WELLS

	Well ID	Hydrostratig raphic Unit Screened	Measuring Point Elevation (feet MSL)	MARCH 2006 Depth to Water (feet)	JUNE 2006 Depth to Water (feet)	Date	Time	Current Well Conditions	
1	MW-1	В		Dry	Δ	R	Y	257'-70 Dry	
2	MW-2	X/A(wt)		Dry	DRY			265'-TD Dry	\Box
3	MW-3	X/A(wt)		244.3	241.56	6/2/16	1230		
4	MW-4	X(wt)		230.6	227.61	6/6/06			\neg
5	MW-5	В		232.28	289-67	416/04	1141		
6	MW-6	A'/X(wt)		230.75	227.96				
7	MW-7	A'/X(wt)		228.3	225.61	6 16 106	1101		
8	MW-8	A'/X(wt)	j.	233.81	227.25		1115		



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TETRA TECH, INC. Fax (628) 351-5291

TC#: 17653-0602

Project: Burbank PAC WELLS

Client: Lockheed Martin Corporation

WELL PURGING FORM Date: 6/20/06

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Sampler NORMAN Monitoring Well ID: Static Water Level (ft bloc): Duplicate ID: 241 Total Well Depth (ft): Well Diameter: 284.50 Water Column (ft): Pump Specs.: 60 TOC to ground surface (ft): Sample Time: TOC = top of casing (at notch/mark) WELL PURGING: 43.00 Notes: Flow RATE - 30 3.4 GPM (water column) 83 4/1+4 (fotal volume to purge) Note: water column x multiplier = casing volume

T12.00		- The state of the								
TIME	(DEG C°)	EC	pH	TURBIDITY and COLOR	SALINITY	DO	TOTAL G			
1821	23.5	0.94	6.88	31	0.04		PURGE			
1124	21,2	0.94	6.87	25	-	6.65				
1129	21.3	0.94	6.86	22	0.04	5,78	10			
1130	21.4	0.94	6.86	20	0.04	5.54	20			
1133	21.3	0.94	6.85	19	0.04	5.50	30			
1136	21.3	0.93	6.85	8	0.04	5.53	40			
1139	21.3	0.94	6.85	4	0.04	5.76	50			
1142	21.3	0.94	6.85	3	0.04	5.,56	60			
1145	21.3	0.94	6.85	3	0.04	5,46	70			
1147	21.3	0.94	6.85	2	0.04	5,42	80			
					0.04	5,45	84			
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Project: Client:

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Burbank PAC WELLS

Lockheed Martin Corporation

Note: water column x multiplier - casing volume

WELL PURGING FORM

Norman Na

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MW-4 Static Water Level (ft bloc): 227 Monitoring Well ID: 264.10 Total Well Depth (ft): Duplicate ID: Water Column (ft): 37 Well Diameter: (No condrol best) TOC to ground surface (ft): Pump Specs.: TOC = top of casing (at notch/mark) Sample Time: WELL PURGING: –4.5 GPM 37.14 Notes: inake (no. of valumes to purge) (lotal volume to purge) 2451

Sampler

TIME	TEMP	EC	рН	TURBIDITY and COLOR	SALINITY	DO	TOTAL GAL.
\$ 309	(DEG C°)	0,93	6.94	195	0,04	5.09	PURGED
1311	20.8	0.92	6.93	165	0.04	4.29	9
1313	20.9	0.742	6.93	34	0.03	4.73	18
1315	21.0	0.742	6.92	24	0.03	449	27
1317	21.0	0.783	6.92	19	0-03	4.36	36
1319	21:0	6.747	691	19	0.03	<i>5</i> 703	45
1321	210	0.746	6.91	7	0.03	4.69	54
1323	21.0	0.748	6.92	5	0-03	4.71	63
1325	21.0	0.746	6.90	7	0.03	4.37	72
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Novmen Na Sampler

Date: _ Page

WELL PURGING FORM

Monitoring Well ID: MW - 5 Duplicate ID: Total Well Depth (ft): 269.50 Well Diameter: 4" Pump Specs.: 3/4 hp 230 V (No control box) Sample Time: 1205 ToC to ground surface (ft): TOC to ground surface (ft): TOC etop of casing (at notch/mark) WELL PURGING: 40.40 x 0.645 = 26 gals (total volume) Well purple of casing volume) ToC to ground surface (ft): TOC etop of casing (at notch/mark) Well purple of casing (at notch/mark) Well purple of casing (at notch/mark) ToC etop of casing (at notch/mark)								
TIME	TEMP (DEG C°)	EC	рН	TURBIDITY and COLOR	SALINITY	DO	TOTAL GAL PURGED	
444	23.6	.95	18	6.93	0.04	5-62	8	
1146	22.8	.91	693	16	0.04	5.00	10	
1148	21.5	.92	6.92	17	0.04	4.74	20	
1150	21.0	.743	694	14	9,04	5.17	30	
1152	21.0	.743	6.93	1.4	0.03	4.70	40	
1154	20.9	-745	6.92	8	0.03	5,21	50	
115%	20.8	746	6.92	8	0.03	5.22	60	
1128	20.8	1746	6.92	8	0.03	5.17	70	
1200	20.9	. 74 4	6-93	6	0-03	4.99	81	

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WELL PURGING FORM

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MW-6 Monitoring Well ID: Static Water Level (ft btoc): 227.80 265.00 Total Well Depth (ft): Duplicate ID: 39.20 Well Diameter: Water Column (ft): pro control TOC to ground surface (ft): Pump Specs.: TOC = top of casing (at notch/mark) Sample Time: WELL PURGING: with Gray Brilling. x 0.645 Notes: 37,20 (water column)

Sampler

24 gals x (1 casing volume)

intake at 2451

Norman Na

Note: water column x multiplier -- casing volume TIME TEMP EC pН TURBIDITY SALINITY DO TOTAL GAL and COLOR **PURGED** (DEG C°) 6.82 .97 6-31 0.04 0923 20.7 206 6.84 \$8 0-04 5-40 18 6.86 0927 20 .-.94 16 \$.50 0-04 10 5.27 20.7 .93 6.87 0.04 0929 0.93 8 36 20.8 687 0931 0.04 5.21 0.93 6-87 3 0933 20.8 0-04 45 <u>5</u>.17 0.92 0935 5.14 54 208 0.04 6-88 6.88 09:37 0.92 20-8 Ó 5.14 0-04 0.92 20.8 6939 6.88 5.15 72 0.04



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Project:

Burbank PAC WELLS

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WELL PURGING FORM

Page

Norman Na Sampler

Monitoring Well ID:	MW-7		Static Water Level (ft btoc):	225.6	
Duplicate ID:			Total Well Depth (ft):	260.0	
Well Diameter:	4"		Water Column (ft):	34.4	
Pump Specs. :	3/4 hp /230V	(no during)	TOC to ground surface (ft):		
Sample Time:	1018		TOC = top of casing (at notch/mark)		
WELL PURGING: 34.4 x	0.645	= 22.2 gals	Notes:	3600	
(water column)	(multiplier)	(1 casing volume)	FLOW RATE -	300 FW	
22.2 gals x		= 66.6 gals	intake at a	250'	
(1 casing volume)	(no. of volumes to purge)	(total volume to purge)			

	multiplier - casing vo						
TIME	TEMP (DEG C°)	EC	рН	TURBIDITY and COLOR	SALINITY	DO	TOTAL GA
0949	21.5	6.91	6.91	128	0.04	6.46	8
0952	21.0	0.738	6.86	120	0.03	5.78	10
0955	21.1	.735	6.88	∂ 8	0.03	5.73	20
0958	21.1	·735	6.88	76	0.63	5.71	30
1001	21.1	.740	6.88	64	0.03	5.81	40
1004	21.1	-738	6.87	22	0.03	5.78	50
1007	21.1	.737 :	6.87	10	0.03	5.86	60
1010	21.1	.738	6.87	6	0.04	5.73	7063
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3475 E. Footh# Blvd. Pasadena, CA 91107 (626) 351-4664

MW-8

TETRA TECH, INC. Pax (628) 351-5291

Monitoring Well ID:

TC#:	17653-0602
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Burbank PAC WELL Project: Client: **Lockheed Martin Corporation**

Sampler

Static Water Level (It bloc):

WELL PURGING FORM Date: 6/20/06 Page

Duplicate ID: Well Diameter: Pump Specs. : Sample Time;	3/4 h p 08	1230 vol	H (no rug		Well Depth (ft): Column (ft): o ground surface top of casing (at	e (ft):	3450 3450		
WELL PURGIN 3 4.20 (water column)	G: x <u>0.64</u> (multiplier)	<u> 5</u>	22 (1 casing volume)	gals Notes	: Flow	RATE -	3 G PM		
22 gals (1 casing volume)	(no. of volume	vols ==	66 (total volume to pu	gels	intale	a at 20	55"		
TIME	TEMP (DEG C°)	EC	pH	TURBIDITY and COLOR	SALINITY	DO	TOTAL GAL PURGED		
0801	20.7	.728	6.76	183	0.04	7.78	0		
0804	20.4	.734	.6.82	157	0.03	5.49	9		
0807	20.8	.732	6.83	28	0.03	5.35	18		
0800	21.0	.733	6.84	20	0.03	5.31	27		
0813	21.1	.732	6.85	12	0.03	5.46	36		
0816	21.3	:733	6.85	6	ტახპ	5.49	45		
0819	21.3	.734	6.85	.4	0.03	5.49	54		
0822	21.3	.734	6.85	4	0.03	5.47	63		
0823	21.3	.734	6.85	4	0.03	5.48	66		
	•								

APPENDIX B LAB ANALYTICAL DATA REPORTS

APPENDIX B

LABORATORY ANALYTICAL DATA REPORTS





July 03, 2006

Neil Shukla Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Subject:

Calscience Work Order No.:

Client Reference:

06-06-1183

BOU Groundwater Monitoring 2006 (PAC Wells)

/ 17653-0602

Dear Client:

Enclosed is an analytical report for the above-referenced project. The samples included in this report were received 6/19/2006 and analyzed in accordance with the attached chain-of-custody.

Unless otherwise noted, all analytical testing was accomplished in accordance with the guidelines established in our Quality Systems Manual, applicable standard operating procedures, and other related documentation. The original report of any subcontracted analysis is provided herein, and follows the standard Calscience data package. The results in this analytical report are limited to the samples tested and any reproduction thereof must be made in its entirety.

If you have any questions regarding this report, please do not hesitate to contact the undersigned.

Sincerely,

Calscience Environmental

Laboratories, Inc.

Jason Torres Project Manager

A-ELAP ID: 1230

NELAP ID: 03220CA

CSDLAC ID: 10109

SCAQMD ID: 93LA0830

7440 Lincoln Way, Garden Grove, CA 92841-1427 • TEL:(714) 895-5494 •

FAX: (714) 894-7501





Tetra Tech, Inc.

Date Received:

06/19/06

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Work Order No:

06-06-1183

Preparation:

EPA 3005A Filt. / EPA 7470A Filt.

Method:

EPA 6010B / EPA 7470A

Units:

mg/L

Client Sample Nu	mber			ample nber		Date Collected	N. A. malania a	ate Da pared Analy	000	atch ID	
MW-6			06-0	6-1183-2		06/19/06	Aqueous 06/	20/06 06/2	21/06 06062	OLO3F	
Comment(s):	-Results were eval	luated to the I	MDL, concentra	tions >= to	the N	MDL but < RL, if for	und, are qualified v	vith a "J" flag.			
	-Mercury was anal	lyzed on 6/20	/2006 3:47:08 P	M with bat	ch 06	0620L02F					
Parameter	Result	<u>RL</u>	MDL.	DF C	Qual	<u>Parameter</u>	Result	<u>RL</u>	MDL	<u>DF</u>	Qua
Antimony	ND	0.0150	0.00209	1		Mercury	ND	0.000500	0.0000672	1	
Arsenic	ND	0.0100	0.00308	1		Molybdenum	ND	0.00500	0.000800	1	
Barium	0.143	0.010	0.000719	1		Nickel	ND	0.00500	0.00137	1	
Beryllium	ND	0.00100	0.000176	1		Selenium	0.00375	0.0150	0.00295	1	J,E
Cadmium	0.00122	0.00500	0.000350	1	J	Silver	ND	0.00500	0.000400	1	
Chromium	0.00793	0.00500	0.000350	1		Thallium	0.0112	0.0150	0.00233	1	J,I
Cobalt	ND	0.00500	0.000696	1		Vanadium	0.00431	0.00500	0.000314	1	,
Copper	ND	0.00500	0.00134	1		Zinc	0.181	0.010	0.000848	1	
ead	ND	0.0100	0.00236	1							
MW-5			06-0	6-1183-3	7.0	06/19/06	Aqueous 06/	20/06 - 66/2	21/06 06062	OLO3F	

<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>MDL</u>	<u>DF</u>	Qual	<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>MDL</u>	<u>DF</u>	Qual
Antimony	ND	0.0150	0.00209	1		Mercury	ND	0.000500	0.0000672	1	
Arsenic	,ND	0.0100	0.00308	1		Molybdenum	ND	0.00500	0.000800	1	
Barium	0.135	0.010	0.000719	1		Nickel	ND	0.00500	0.00137	1	
Beryllium	ND	0.00100	0.000176	1		Selenium	0.00393	0.0150	0.00295	1	J,B
Cadmium	ND	0.00500	0.000350	1		Silver	ND	0.00500	0.000400	1	
Chromium	0.00648	0.00500	0.000350	1		Thailium	0.00498	0.0150	0.00233	1	J,B
Cobalt	· ND	0.00500	0.000696	1		Vanadium	0.00353	0.00500	0.000314	. 1	J
Copper	ND	0.00500	0.00134	1		Zinc	0.144	0.010	0.000848	1	
Lead	ND	0.0100	0.00236	1							

MW-4 06-06-1183-4 06/19/06 Aqueous 06/20/06 06/21/06 060620L03F

Comment(s): -Results were evaluated to the MDL, concentrations >= to the MDL but < RL, if found, are qualified with a "J" flag.

-Mercury was analyzed on 6/20/2006 3:51:36 PM with batch 060620L02F

<u>Parameter</u>	Result	<u>RL</u>	<u>MDL</u>	DF C	<u>Qual Parameter</u>	Result	<u>RL</u>	MDL	DF Qual
Antimony	ND	0.0150	0.00209	1	Mercury	ND	0.000500	0.0000672	1
Arsenic	ND	0.0100	0.00308	1	Molybdenum	ND	0.00500	0.000800	1
Barium	0.146	0.010	0.000719	1	Nickel	ND	0.00500	0.00137	1
Beryllium	ND	0.00100	0.000176	1	Selenium	ND	0.0150	0.00295	1
Cadmium	ND	0.00500	0.000350	1	Silver	ND	0.00500	0.000400	1
Chromium	0.00561	0.00500	0.000350	1	Thallium	0.00694	0.0150	0.00233	1 ј
Cobalt	ND	0.00500	0.000696	1	Vanadium	0.00477	0.00500	0.000314	1 ј
Copper	ND	0.00500	0.00134	1	Zinc	0.0850	0.0100	0.000848	1
Lead	ND	0.0100	0.00236	1					

RL - Reporting Limit ,

DF - Dilution Factor ,







Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

06/19/06

Work Order No:

06-06-1183

Preparation:

Method:

EPA 3005A Filt. / EPA 7470A Filt.

Date

Date

Date

EPA 6010B / EPA 7470A

Units:

mg/L

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Page 2 of 2

Client Sample Nu	mber		Nur	nber	Collected	Matrix Pre	pared Anal	yzed QCB	atch ID	
MW-4D			06-0	6-1183-5	06/19/06	Aqueous 06	20/06 🕮 06/	21/06 06062	OLO3F	, E
Comment(s):	-Results were eva	luated to the I	MDL, concentrat	tions >= to the	MDL but < RL, if f	found, are qualified	with a "J" flag.			
	-Mercury was ana	lyzed on 6/20	/2006 3:58:21 P	M with batch (60620L02F					
<u>Parameter</u>	Result	<u>RL</u>	MDL	<u>DF</u> Qua	<u>Parameter</u>	Result	<u>RL</u>	MDL	DF !	Qual
Antimony	ND	0.0150	0.00209	1	Mercury	ND	0.000500	0.0000672	1	
Arsenic	ND	0.0100	0.00308	1	Molybdenum	ND	0.00500	0.000800	1	
Barium	0.145	0.010	0.000719	1	Nickel	ND	0.00500	0.00137	1	
Beryllium	ND	0.00100	0.000176	1	Selenium	0.00831	0.0150	0.00295	1	J,B
Cadmium	ND	0.00500	0.000350	1 1	Silver	ND	0.00500	0.000400	1	
Chromium	0.00622	0.00500	0.000350	1	Thailium	0.00594	0.0150	0.00233	1	J,B
Cobalt	ND	0.00500	0.000696	1	Vanadium	0.00467	0.00500	0.000314	1	J
Copper	ND	0.00500	0.00134	1	Zinc	0.122	0.010	0.000848	1	
Lead	ND	0.0100	0.00236	1						

Method Blank 099-04-008-2,541 N/A Aqueous 06/20/06 06/20/06 060620L02F

-Results were evaluated to the MDL, concentrations >= to the MDL but < RL, if found, are qualified with a "J" flag.

Lab Sample

<u>Parameter</u> Mercury

Result ND

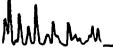
0.000500

DF Qual

Method Blank			097-	01-003-6	215	NA A	queous 06/20)/06 - 06/	21/06 06062	0L03F	
Comment(s):	-Results were eva	luated to the I	MDL, concentra	tions >=	to the i	MDL but < RL, if four	nd, are qualified wit	th a "J" flag.			
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	MDL	<u>DF</u>	<u>Qual</u>	<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>MDL</u>	DF 9	Qual
Antimony	ND	0.0150	0.00209	1		Lead	ND	0.0100	0.00236	1	
Arsenic	0.00391	0.0100	0.00308	1	J	Molybdenum	ND	0.00500	0.000800	1	
Barium	ND	0.0100	0.000719	1		Nickel	0.00394	0.00500	0.00137	1	J
Beryllium	ND	0.00100	0.000176	1		Selenium	0.00502	0.0150	0.00295	1	J
Cadmium	ND	0.00500	0.000350	1		Silver	ND	0.00500	0.000400	1	
Chromium	ND	0.00500	0.000350	1		Thallium	0.00431	0.0150	0.00233	1	J
Cobalt	ND	0.00500	0.000696	1		Vanadium	ND	0.00500	0.000314	1	
Copper	ND	0.00500	0.00134	1		Zinc	ND	0.0100	0.000848	1	



DF - Dilution Factor







Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

Work Order No:

Preparation:

Method: Units:

06/19/06

06-06-1183 EPA 3005A Filt.

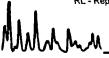
EPA 6010B

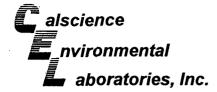
mg/L

Project: BO	U Groundwat	ter Monit	oring 2006 (PAC V	Vells) / 17653-0	0602			Page 1 of	<u>1</u>
Client Sample Nu	mber			Sample mber		Date Collected	Matrix	Date Prepared	Date Analyzed	QC Batch ID	
MW-6 2 4 2 2		7 Y	w> 06-0	6-1183-2		1 06/19/06	Aqueous	06/20/06	06/21/06	060620L03F	
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	MDL	<u>DF</u>	Qual	<u>Parameter</u>	Res	ult RL	<u>MDL</u>	<u>D</u> F	Qual
Calcium	112	0.100	0.00932	1		Potassium		5.87 0.5	0.056	1 1	l
Magnesium	30.8	0.1	0.00328	1	В	Sodium		37.0 0.5	0.019	2 1	
MW-5			06-0	6-1183-3		06/19/06	Aqueous	06/20/06	06/21/06	060620L03F	
Parameter	Result	RL	MDL	DE	Qual	<u>Parameter</u>	Res	sult RL	MDL	<u>DE</u>	Qual
Calcium	109	0.100	0.00932	1		Potassium		5.66 0.5	0 0.056	- 11 1	1
Magnesium	29.9	0.1	0.00328	1	В	Sodium	: :	36.60.5	0.019	2 1	l
MW-4			. 06-0	6-1183-4		06/19/06	Aqueous	08/20/06	06/21/06	060620L03F	
<u>-</u>										_	
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>MDL</u>	<u>DF</u>	Qual	<u>Parameter</u>	Res	sult RL	<u>MDL</u>	<u>D</u> F	Qual
Calcium	104	0.100	0.00932	1	Ε	Potassium		5.56 0.5	0 0.056	31 1	1
Magnesium	33.7	0.1	0.00328	1		Sodium		37.9 0.5	0.019)2 1	<u> </u>
MW-4D			06-0	6-1183-5		06/19/06	Aqueous	06/20/06	06/21/06	060620L03F	
				,							
Parameter Parameter	Result	<u>RL</u>	<u>MDL</u>	DF	Qual	<u>Parameter</u>	Res	sult RL	MDL	DF	Qual
Calcium	104	0.100	0.00932	1		Potassium		5.41 0.5	0 0.056	31 1	I
Magnesium	33.4	. 0.1	0.00328	1	В	Sodium		37.4 0.5	0.019	21	<u> </u>
Method Blank	and the second	动性	097-	01-003-6	,215	NA	Aqueous	06/20/06	06/21/06	060620L03F	
Comment(s):	-Results were eva	aluated to the	e MDL, concentra	itions >=	to the N	MDL but < RL, i	f found, are qu	alified with a "J	" flag.		
<u>Parameter</u>	Result	<u>RL</u>	MDL	<u>DF</u>	Qual	<u>Parameter</u>	Res	sult RL	MDL	<u>DF</u>	Qual
Calcium	ND	0.100	0.00932	1		Potassium		ID 0.5	00 0.056	31 · 1	1
Magnesium	0.0152	0.100	0.00328	1	J	Sodium	N	ID 0.5	0.019)2 1	i

RL - Reporting Limit ,

DF - Dilution Factor







Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

eceived: 06/19/06

Work Order No: 06-06-1183

Preparation: EPA 3005A Filt. Method: EPA 200.8

Units:

mg/L

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Page 1 of 1

			Lab	Sample		Date		Date		Date		
lient Sample Nu	mber		NuNu	ımber		Collected	Matrix	Prepare	ed .	Analyzed	QC Bat	ich ID
MW-6			. 06-	06-1183-2		06/19/06	Aqueous	06/21/0	16	06/21/06	060621	IL01 🙏 🛬
Comment(s):	-Results were eva	aluated to the	MDL, concentr	ations >= 1	to the f	MDL but < RL, if	found, are qua	alified with	a "J" fl	lag.		
Parameter	<u>Result</u>	<u>RL</u>	<u>MDL</u>	<u>DF</u>	<u>Qual</u>	<u>Parameter</u>	Res		<u>RL</u>	<u>MDL</u>	•	DF Qu
ron	0.0260	0.100	0.00214	1	J	Manganese	0.	00439	0.001	00.000	0189	1
MW-5 🚁 🐰			.06-	06-1183-3		22. 06/19/06	Aqueous	06/21/0)6	06/21/06	060621	L01 🐪
Comment(s):	-Results were eva	aluated to the	MDL, concentr	ations >= 1	to the f	MDL but < RL, if	found, are qua	alified with	a "J" fl	lag.		
arameter	Result	<u>RL</u>	MDL	<u>DF</u>	Qual	<u>Parameter</u>	Res	<u>ult</u>	<u>RL</u>	MDL	•	<u>DF</u> Qu
on	0.0301	0.100	0.00214	1	J	Manganese	0.	00261	0.001	0.000	0189	, 1
				Market British and Carlo Control of the	Bur Est and	APPA DA BASANDON CONTRARADOR DE CONTRARADOR DE CONTRARADOR DE CONTRARADOR DE CONTRARADOR DE CONTRARADOR DE CONT	Mary Section 12 13 Per	\$3.00 F	TO SHOW	Calabra Mar Principle	COMMENCE:	第4个公司和
MW-4			# 2 ¹ = ₹ 06-	06-1183-4		06/19/06	Aqueous :	06/21/0)6	06/21/06	060621	ILO1
MW-4 Comment(s):	-Results were eva	aluated to the	No. Address Professor Start of the supplies of	er and and the second	SECTION AND ADDRESS.	# 198 market 12 \$ 20-11 def-140000 1000000	Market Cont	· · · · · · · · · · · · · · · · · · ·	(1) - COMMO	enin adampte (allahara)	060621	ILO1
Comment(s):	-Results were eva	aluated to the RL	No. Address Professor Start of the supplies of	ations >=	SECTION AND ADDRESS.	MDL but < RL, if	Market Cont	alified with	(1) - COMMO	enin adampte (allahara)	<u> </u>	DF Qu
Comment(s):			MDL, concentr	ations >=	to the f	MDL but < RL, if	found, are qua	alified with	a "J" fl	lag.		
Comment(s):	Result	<u>RL</u>	MDL, concentr MDL 0.00214	ations >=	to the I	MDL but < RL, if Parameter	found, are qua	alified with ult 00497	a "J" fl <u>RL</u> 0.001	lag.		<u>DF Qu</u> 1
Comment(s): <u>Parameter</u> on	Result	RL 0.100	MDL, concentr MDL 0.00214	ations >= 1 <u>DF</u> 1 06-1183-5	to the f Qual J	MDL but < RL, if Parameter Manganese 06/19/06	found, are qua Res 0. Aqueous	alified with ult 00497 06/217	a "J" fl RL 0.001	lag. MDL 00 0.000	00189	<u>DF Qu</u> 1
Comment(s): Parameter Ton MW4D Comment(s):	Result 0.0342	RL 0.100	MDL, concentr MDL 0.00214	ations >= 1 <u>DF</u> 1 06:1183-5 ations >= 1	to the I Qual J	MDL but < RL, if Parameter Manganese 06/19/06	found, are qua Res 0. Aqueous	alified with ult 00497 06/217	a "J" fl RL 0.001	lag. MDL 00 0.000	00189 06062:	<u>DF Qu</u> 1
Comment(s): Parameter ron Comment(s): Parameter	Result 0.0342	RL 0.100	MDL, concentr MDL 0.00214	ations >= 1 <u>DF</u> 1 06:1183-5 ations >= 1	to the I Qual J	MDL but < RL, if Parameter Manganese 06/19/06	found, are quantum found, are qu	alified with ult 00497 06/217	a "J" fl <u>RL</u> 0.001 06	lag. MDL 00 0.000 06/21/06	- 00189 - 06062 *	DF Qu 1
Comment(s): Parameter con Comment(s): Parameter	Result 0.0342 -Results were eva Result 0.0246	RL 0.100 aluated to the	MDL, concentr MDL 0.00214 0.00214 MDL, concentr MDL 0.00214	ations >= 1 <u>DF</u> 1 06:1183-5 ations >= 1	to the I Qual J to the I Qual J	MDL but < RL, if Parameter Manganese 06/19/06 MDL but < RL, if Parameter	found, are quantum found, are qu	alified with ult 00497 06/21# alified with ult 00488	a "J" fi RL 0.001 0.001 a "J" fi RL 0.001	lag. MDL 00 0.000 06/21/06	- 00189 - 06062 *	DF Qu 1 ILO1 DF Qu 1
Comment(s): Parameter on MW-4D Comment(s): Parameter on	Result 0.0342 -Results were eva Result 0.0246	RL 0.100 aluated to the RL 0.100	MDL, concentr MDL 0.00214 06- MDL, concentr MDL 0.00214	ations >= 1 DF 1 06-1183-5 ations >= 1 DE 1 1-10-008-7	to the f Qual J to the f Qual J	MDL but < RL, if Parameter Manganese 06/19/06 MDL but < RL, if Parameter Manganese	found, are quantities of the country	alified with ult 00497 06/21R alified with ult 00488	a "J" fi RL 0.001 06 a "J" fi RL 0.001	lag. 00 0.000 06/21/06 lag. MDL 00 0.000 06/21/06	00189	DF Qu 1 ILO1 DF Qu 1
Comment(s): Parameter ron Comment(s): Parameter ron Method Blank	Result 0.0342 -Results were ever Result 0.0246	RL 0.100 aluated to the RL 0.100	MDL, concentr MDL 0.00214 06- MDL, concentr MDL 0.00214	ations >= 1 DF 1 06-1183-5 ations >= 1 P-10-008-7 ations >= 1	to the f Qual J to the f Qual J	MDL but < RL, if Parameter Manganese 06/19/06 MDL but < RL, if Parameter Manganese N/A	found, are quantities of the country	alified with ult 00497 06/21R alified with ult 00488 06/21/c	a "J" fi RL 0.001 06 a "J" fi RL 0.001	lag. 00 0.000 06/21/06 lag. MDL 00 0.000 06/21/06	00189 06062* 00189	DF Qu 1 ILO1 DF Qu 1

RL - Reporting Limit

DF - Dilution Factor





Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

Work Order No:

Preparation: Method:

06-06-1183

EPA 3520B

EPA 8270C(M) Isotope

Dilution

06/19/06

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Page 1 of 2

Client Sample Number		Lab Samp Number		Date Collected	Matrix	Date Prepared	Date Analyzed	QC Batch ID
MVV-8		06,06-1	83-2: 🔭	\$, 06/19/06	Aqueous	06/22/06	06/28/06	060622L01D
Comment(s): -Results were ev	aluated to the	MDL, concentrations	s >= to the M	DL but < RL, if	found, are qua	lified with a "J"	' flag.	-
<u>Parameter</u>	Result	<u>RL</u>	MDL	<u>DF</u>	<u>Qual</u>	<u>Units</u>		
1,4-Dioxane	ND	2.0	0.40	1		ug/L		
Surrogates:	<u>REC (%)</u>	Control Limits			<u>Qual</u>			
Nitrobenzene-d5	80	56-123						<u> </u>
MW-5 (1)		The state of the s	83-3/4 - ***		Aqueous	06/22/06	06/28/06	:: 060622L01D
Comment(s): -Results were ev		•			•		'flag.	
<u>Parameter</u>	Result	<u>RL</u>	<u>MDL</u>	DE	<u>Qual</u>	<u>Units</u>		
1,4-Dioxane	ND	2.0	0.40	. 1		ug/L		
Surrogates:	<u>REC (%)</u>	Control Limits			<u>Qual</u>			
Nitrobenzene-d5	82	56-123						
MW-4		* 06-06-1	183-4	06/19/06	Aqueous	06/22/06	06/28/06	060622L01D
Comment(s): -Results were ev					_		' flag.	
<u>Parameter</u>	Result	<u>RL</u>	<u>MDL</u>	<u>DF</u>	<u>Qual</u>	<u>Units</u>		
1,4-Dioxane	ND	2.0	0.40	1		ug/L		
Surrogates:	<u>REC (%)</u>	Control Limits			<u>Qual</u>			
Nitrobenzene-d5	78	56-123				•		-
MW-4D		06-06-1	183-5	.06/19/06	Aqueous	06/22/06	06/28/06	060622L01D
Comment(s): -Results were ev		·		· ·	•		' flag.	
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>MDL</u>	<u>DF</u>	<u>Qual</u>	<u>Units</u>		
1,4-Dioxane	ND	2.0	0.40	1		ug/L		
Surrogates:	<u>REC (%)</u>	Control Limits			Qual			
Nitrobenzene-d5	67	56-123						
Method Blank			04-596	NA	Aqueous	06/20/06	06/27/06	060620L14D
<u>Parameter</u>	Result	<u>RL</u>	MDL	<u>DF</u>	<u>Qual</u>	<u>Units</u>		
1,4-Dioxane	ND	2.0		1		ug/L		
Surrogates:	REC (%)	Control Limits		•	<u>Qual</u>	-g, -		
Nitrobenzene-d5	07	56-123		•				
INIU ODENZENE-OD	87	JU-123						

RL - Reporting Limit

DF - Dilution Factor







Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

Work Order No:

Preparation:

Method:

06/19/06

06-06-1183

EPA 3520B

EPA 8270C(M) Isotope

Dilution

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Page 2 of 2

Client Sample Number		Lab Sam Numbe		Date Collected	Matrix	Date Prepared	Date Analyzed	QC Batch ID
Method Blank	4	099-09-0	004-597,	₽ NA	Aqueous	06/22/06	06/27/06	060622L01D
Comment(s): -Results were eval	uated to the	MDL, concentrations	s >= to the MI	DL but < RL, if	found, are qua	lified with a "J"	flag.	
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	MDL	<u>DF</u>	<u>Qual</u>	<u>Units</u>		
1,4-Dioxane	ND	2.0	0.40	1		ug/L		
Surrogates:	REC (%)	Control Limits			<u>Qual</u>			
Nitrobenzene-d5	76	56-123						







Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

Work Order No:

Preparation: Method:

06-06-1183 **EPA 3520B**

06/19/06

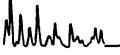
EPA 1625CM

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Page 1 of 1

Client Sample Number		Lab Sampl Number	e	Date Collected	Matrix	Date Prepared	Date Analyzed	QC Batch ID
MW-6		06-06-118	3-2	06/19/06	Aqueous	06/23/06	06/27/06	060623L09
Comment(s): -Results were eva							flag.	
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	MDL	DF	<u>Qual</u>	<u>Units</u>		
N-Nitrosodimethylamine Surrogates:	ND REC (%)	2.0 Control Limits	0.48	1	Qual	ng/L		
1,4-Dichlorobenzene-d4	56	50-130				_		
MW-5		₹06-06-118	MAN PARTICIPATION OF THE	The state of the s	Aqueous	06/23/06	** 06/27/06 ,	060623L09 4
Comment(s): -Results were eva					-		flag.	
<u>Parameter</u>	Result	<u>RL</u>	MDL	<u>DF</u>	<u>Qual</u>	<u>Units</u>		
N-Nitrosodimethylamine	ND	2.0	0.48	1		ng/L		
Surrogates:	REC (%)	Control Limits			<u>Qual</u>			
1,4-Dichlorobenzene-d4	53	50-130						
MW-4		06-06-111	83-4 , * (06/19/06	Aqueous	² 06/23/06	06/27/06	060623L09
Comment(s): -Results were eva							flag.	
<u>Parameter</u>	Result	<u>RL</u>	MDL	<u>DF</u>	<u>Qual</u>	<u>Units</u>		
N-Nitrosodimethylamine	ND	2.0	0.48	1		ng/L		
Surrogates:	REC (%)	Control Limits			<u>Qual</u>			
1,4-Dichlorobenzene-d4	55	50-130	_					
MW-ID:		06-06-118	100000000000000000000000000000000000000	06/19/06	Aqueous	06/23/06	≢ 06/27/06	060623L09
Comment(s): -Results were eva							flag.	
Parameter	Result	RL	<u>MDL</u>	<u>DF</u>	Qual	<u>Units</u>		
N-Nitrosodimethylamine	ND	2.0	0.48	1		ng/L		
Surrogates:	REC (%)	Control Limits			<u>Qual</u>			
1,4-Dichlorobenzene-d4	59	50-130						
Method Blank		099-07-0	ent Spirothic and Co.	. NA	Aqueous	06/23/06	06/27/06	060623L09:
Comment(s): -Results were eva Parameter	lluated to the Result	MDL, concentrations <u>RL</u>	>= to the M MDL	DL but < RL, if DF	found, are qua Quai	alified with a "J" <u>Units</u>	flag.	
N-Nitrosodimethylamine	ND	2.0	0.48	1	Ovel	ng/L		
Surrogates:	REC (%)	Control Limits			<u>Qual</u>			
1,4-Dichlorobenzene-d4	54	50-130						

DF - Dilution Factor







Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

06/19/06

Work Order No:

06-06-1183 **EPA 5030B**

Preparation:

Method:

EPA 8260B

Units:

ug/L

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Page 1 of 8

Comment(s): Results were evaluated to the MDL, concentrations >= to the MDL but < RL, if found, are qualified with a "J" flag.	Client Sample Number			Lab Sa Numi			Date Collected	Matrix	Date Prepared	Date Analyzed	QC Bate	ch ID	
Parameter Result RI. MDI. DE Qual Parameter Result RI. MDI. DE Qual Acetone 6.4 10.0 6.1 1 J. 3.3 - 1.3	LTB061906			06-06-	1,183-1		06/19/06	Aqueous	06/22/06	06/23/06	060622	L02	. 4
Acetone 6.4 10.0 6.1 1 J.B 1,3-Dichloropropane ND 1,0 0.30 1 Bervene ND 0.50 0.26 1 2,2-Dichloropropane ND 1,0 0.40 1 Bromochloromethane ND 1,0 0.68 1 c-1,3-Dichloropropene ND 0.50 0.45 1 Bromochloromethane ND 1,0 0.68 1 c-1,3-Dichloropropene ND 0.50 0.45 1 Bromofform ND 1,0 0.62 1 Ethylchichoropropene ND 0.50 0.45 1 Bromofform ND 1,0 0.62 1 Ethylchichoropropene ND 0.0 0.01 1 Bromofform ND 1,0 0.62 1 Ethylchichoropropene ND 0.0 0.01 0.0 0.01 0.0 0.01 0.00 0.00 0.00 0.00 0.00 0.00 0.00 0.00 0.00 <	Comment(s): -Results were	evaluated to the	e MDL, co	ncentratio	ons >= t	o the M	MDL but < RL, if f	ound, are qualifi	ed with a "J" f	ag.			
Benzene ND 0.50 0.26 1 2,2-Dichloropropane ND 1.0 0.40 1	<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>MDL</u>	<u>DF</u>	Qual	<u>Parameter</u>		Result	<u>RL</u>	<u>MDL</u>	<u>DF</u> Q	<u>ual</u>
Bromobenzene ND 1.0 0.47 1 1,1-Dichloropropene ND 1.0 0.21 1 Bromochloromethane ND 1.0 0.68 1 -1,3-Dichloropropene ND 0.50 0.45 1 Bromochloromethane ND 1.0 0.27 1 +1,3-Dichloropropene ND 0.50 0.45 1 Bromodichoromethane ND 1.0 0.27 1 +1,3-Dichloropropene ND 0.50 0.31 1 Bromodromethane ND 1.0 0.62 1 Ethylbenzene ND 1.0 0.17 1 Bromomethane ND 10 2.9 1 2-Hexanone ND 10 1.9 1 Bromodichoromethane ND 1.0 0.29 1 Stopropythenzene ND 1.0 0.24 1 Bromodichoromethane ND 1.0 0.29 1 Stopropythenzene ND 1.0 0.24 1 Bromodichoromethane ND 1.0 0.21 1 Methylene Chloride 6.1 10.0 0.2 1 Bromodichoromethane ND 1.0 0.17 1 4-Methyl-2-Pentanone ND 1.0 0.2 1 Bromodichoromethane ND 1.0 0.35 1 Naphthalene ND 1.0 0.55 1 Bromodichoromethane ND 1.0 0.52 1 Naphthalene ND 1.0 0.30 1 Bromodichoromethane ND 1.0 0.52 1 Naphthalene ND 1.0 0.30 1 Bromodichoromethane ND 1.0 0.22 1 Naphthalene ND 1.0 0.30 1 Bromodichoromethane ND 1.0 0.32 1 Naphthalene ND 1.0 0.39 1 Bromodichoromethane ND 1.0 0.32 1 Naphthalene ND 1.0 0.39 1 Bromodichoromethane ND 1.0 0.32 1 Naphthalene ND 1.0 0.39 1 Bromodichoromethane ND 1.0 0.32 1 Naphthalene ND 1.0 0.33 1 Bromodichoromethane ND 1.0 0.32 1 Naphthalene ND 1.0 0.33 1 Bromodichoromethane ND 1.0 0.32 1 Naphthalene ND 1.0 0.33 1 Bromodichoromethane ND 1.0 0.32 1 Naphthalene ND 1.0 0.33 1 Bromodichoromethane ND 1.0 0.32 1 Naphthalene ND 1.0 0.33 1 Bromodichoromethane ND 1.0 0.33 1 Naphthalene ND 1.0 0.33 1 Bromodichoromethane ND 1.0 0.33 1 Naphthalene ND 1.0 0.33 1 Bromodichoromethane ND 1.0 0.34 1 Naphthalene ND 1.0 0.35 1 Bro	Acetone	6.4	10.0	6.1	1	J,B	1,3-Dichloropro	pane	ND	1.0	0.30	1	
Bromochtoromethane ND 1.0 0.47 1 1.1-Dichtoropropene ND 0.50 0.21 1	Benzene	ND	0.50	0.26	1		2.2-Dichloropro	pane	ND	1.0	0.40	1	
Bromodichloromethane ND 1.0 0.68 1 c-1,3-Dichloropropene ND 0.50 0.45 1	Bromobenzene	ND	1.0	0.47	1				ND	1.0	0.21	1	
Bromodichloromethane ND 1.0 0.27 1 Ethylbenzene ND 0.50 0.31 1 1 1 1 1 1 1 1 1	Bromochloromethane	ND	1.0	0.68	1				ND	0.50	0.45	1	
Bromomethane	Bromodichloromethane	ND	1.0	0.27	1		•	•	ND	0.50	0.31	1	•
2-Butanone ND 10 4.2 1	Bromoform	ND	1.0	0.62	1			•	ND	1.0	0.17	1	
n-Butylbenzene ND 1.0 0.28 1 p-Isopropyltoluene ND 1.0 0.21 1 sec-Butylbenzene ND 1.0 0.21 1 Methylene Chloride 6.1 10.0 2.6 1 J,B tert-Butylbenzene ND 1.0 0.17 1 4-Methyl-2-Pentanone ND 10 2.4 1 Carbon Disulfide ND 1.0 0.50 0.42 1 n-Propylbenzene ND 1.0 0.95 1 Carbon Tetrachloride ND 1.0 0.36 1 Styrene ND 1.0 0.29 1 Chlorothane ND 1.0 0.52 1 1,1,1,2-Tetrachloroethane ND 1.0 0.29 1 Chlorothane ND 1.0 0.22 1 1,1,2,2-Tetrachloroethane ND 1.0 0.37 1 Chlorothane ND 1.0 0.24 1 Toluene ND 1.0 0.35 1	Bromomethane	ND	10	2.9	1		2-Hexanone		ND	10	1.9	1	
Sec-Butylbenzene ND 1.0 0.21 1 Methylene Chloride 6.1 10.0 2.5 1 J,B	2-Butanone	ND	10	4.2	1		Isopropylbenze	ne	ND	1.0	0.24	1	
tert-Butylbenzene ND 1.0 0.17 1 4-Methyl-2-Pentanone ND 10 2.4 1 Carbon Disulfide ND 10 1.0 1.0 1.0 1.0 1.0 0.95 1 Carbon Tetrachloride ND 0.50 0.42 1 n-Propylbenzene ND 1.0 0.30 1 Chlorobenzene ND 1.0 0.35 1 Styrene ND 1.0 0.29 1 Chloroform ND 1.0 0.52 1 1,1,1,2-Tetrachloroethane ND 1.0 0.37 1 Chloroform ND 1.0 0.52 1 1,1,1,2-Tetrachloroethane ND 1.0 0.37 1 Chlorofoluene ND 1.0 0.22 1 1,1,2-Tetrachloroethane ND 1.0 0.35 1 2-Chlorotoluene ND 1.0 0.24 1 Totuene ND 1.0 0.39 1 1-Chlorotoluene <th< td=""><td>n-Butylbenzene</td><td>ND</td><td>1.0</td><td>0.29</td><td>1</td><td></td><td>p-Isopropyltolue</td><td>ene</td><td>ND</td><td>1.0</td><td>0.21</td><td>1</td><td></td></th<>	n-Butylbenzene	ND	1.0	0.29	1		p-Isopropyltolue	ene	ND	1.0	0.21	1	
Carbon Disulfide ND 10 1.0 1 Naphthalene ND 10 0.95 1 Carbon Tetrachloride ND 0.50 0.42 1 n-Propylbenzene ND 1.0 0.30 1 Chlorobenzene ND 1.0 0.52 1 n-Propylbenzene ND 1.0 0.29 1 Chloroform ND 1.0 0.52 1 1,1,1,2-Tetrachloroethane ND 1.0 0.37 1 Chloroform ND 1.0 0.22 1 1,1,1,2-Tetrachloroethane ND 1.0 0.37 1 Chloroformethane ND 1.0 0.22 1 1,1,2-Tetrachloroethane ND 1.0 0.39 1 2-Chlorotoluene ND 1.0 0.30 1 1,2,3-Trichloroethene ND 1.0 0.35 1 2-Chlorotoluene ND 1.0 0.45 1 1,2,4-Trichloroethane ND 1.0 0.35 1 1	sec-Butylbenzene	ND	1.0	0.21	1		Methylene Chlo	oride	6.1	10.0	2.6	1	J,B
Carbon Tetrachloride ND 0.50 0.42 1 n-Propylbenzene ND 1.0 0.30 1 Chlorobenzene ND 1.0 0.36 1 Styrene ND 1.0 0.29 1 Chlorobenzene ND 1.0 0.52 1 1,1,1,2-Tetrachloroethane ND 1.0 0.37 1 Chloroform ND 1.0 0.22 1 1,1,1,2-Tetrachloroethane ND 1.0 0.37 1 Chlorofoluene ND 1.0 0.22 1 1,1,2-Tetrachloroethane ND 1.0 0.35 1 4-Chlorofoluene ND 1.0 0.24 1 Toluene ND 1.0 0.35 1 4-Chlorotoluene ND 1.0 0.24 1 1,2-Trichloroethane ND 1.0 0.35 1 1/2-Dibromochloromethane ND 1.0 0.45 1 1,2-Trichloroethane ND 1.0 0.32 1 1/2-Di	tert-Butylbenzene	ND	1.0	0.17	1		4-Methyl-2-Pen	itanone	ND	10	2.4	1	
Chlorobenzene ND 1.0 0.36 1 Styrene ND 1.0 0.29 1 Chloroethane ND 1.0 0.52 1 1,1,1,2-Tetrachloroethane ND 1.0 0.37 1 Chloroform ND 1.0 0.22 1 1,1,2,2-Tetrachloroethane ND 1.0 0.37 1 Chloroformethane ND 1.0 0.22 1 1,1,2,2-Tetrachloroethane ND 1.0 0.37 1 2-Chlorotoluene ND 1.0 0.24 1 Toluene ND 1.0 0.35 1 4-Chlorotoluene ND 1.0 0.35 1 1,2,3-Trichlorobenzene ND 1.0 0.35 1 4-Chlorotoluene ND 1.0 0.45 1 1,2,4-Trichlorobenzene ND 1.0 0.39 1 1/2-Dibromo-3-Chloropropane ND 1.0 0.45 1 1,1,1-Trichloroethane ND 1.0 0.32 1	Carbon Disulfide	ND	10	1.0	1		Naphthalene	•	ND	10	0.95	1	
Chloroftane ND 1.0 0.52 1 1,1,1,2-Tetrachloroethane ND 1.0 0.37 1 Chloroform ND 1.0 0.22 1 1,1,1,2-Tetrachloroethane ND 1.0 0.37 1 Chloroftane ND 1.0 0.22 1 1,1,1,2-Tetrachloroethane ND 1.0 0.37 1 Chloroftane ND 1.0 0.24 1 Tetrachloroethane ND 1.0 0.35 1 -Chlorotoluene ND 1.0 0.30 1 1,2,3-Trichlorobenzene ND 1.0 0.35 1 -Chlorotoluene ND 1.0 0.45 1 1,2,4-Trichlorobenzene ND 1.0 0.35 1 -Chlorotoluene ND 1.0 0.45 1 1,2,4-Trichloroethane ND 1.0 0.35 1 -Chlorotoluene ND 1.0 0.45 1 1,1,1-Trichloroethane ND 1.0 0.35 1 -Chlorotoluene ND 1.0 0.45 1 1,1,1-Trichloroethane ND 1.0 0.35 1 -Chlorotoluene ND 1.0 0.45 1 1,1,1-Trichloroethane ND 1.0 0.35 1 -Chlorotoluene ND 1.0 0.45 1 1,1,1-Trichloroethane ND 1.0 0.32 1 -Chlorotoluene ND 1.0 0.81 1 1,1,1-Trichloroethane ND 1.0 0.32 1 -Chlorotoluene ND 1.0 0.81 1 1,1,1-Trichloroethane ND 1.0 0.32 1 -Chlorotoluene ND 1.0 0.42 1 1,1,2-Trichloro-1,2,2-Trifluoroethane ND 1.0 0.54 1 -Chlorotoluene ND 1.0 0.38 1 Trichlorofluoromethane ND 1.0 0.30 1 -Chlorotoluene ND 1.0 0.30 1 1,2,3-Trichloropropane ND 1.0 0.36 1 -Chlorotoluene ND 1.0 0.53 1 1,2,3-Trichloropropane ND 1.0 0.26 1 -Chlorotoluene ND 1.0 0.53 1 1,3,5-Trimethylbenzene ND 1.0 0.30 1 -Chlorotoluene ND 1.0 0.35 1 1,3,5-Trimethylbenzene ND 1.0 0.30 1 -Chlorothane ND 1.0 0.35 1 1,3,5-Trimethylbenzene ND 1.0 0.30 1 -Chlorothane ND 1.0 0.35 1 1,3,5-Trimethylbenzene ND 1.0 0.30 1 -Chlorothane ND 1.0 0.35 1 1,3,5-Trimethylbenzene ND 1.0 0.30 1 -Chlorothane ND 1.0 0.35 1 1,3,5-Trimethylbenzene ND 1.0 0.30 1 -Chlorothane ND 1.0 0.35 1 1,3,5-Trimethylbenzene ND 1.0 0.30 1 -Chlorothane ND 1.0 0.35 1 1,3,5-Trimethylbenzene ND 1.0 0.30 1 -Chlorothane ND 1.0 0.35 1 1,3,5-Trimethylbenzene ND 1.0 0.30 1 -Chlorothane ND 1.0 0.35 1 1,3,5-Trimethylbenzene ND 1.0 0.30 1 -Chlorothane ND 1.0 0.35 1 1,3,5-Trimethylbenzene ND 1.0 0.30 1 -Chlorothane ND 1.0 0.35 1 1,3,5-Trimethylbenzene ND 1.0 0.30 1 -Chlorothane ND 1.0 0.35 1 1,3,5-Trimethylbenzene ND 1.0 0.30 1 -Chlorothane ND 1.0 0.35 1 1,2,4-Trichlorothane ND 1.0 0.30 1 -Chlorothane ND 1.0 0.35 1	Carbon Tetrachloride	ND	0.50	0.42	1		n-Propylbenzer	ne e	ND	1.0	0.30	1	
Chloroform	Chlorobenzene	ND	1.0	0.36	1		Styrene		ND	1.0	0.29	1	
Chloromethane ND 10 1.8 1 Tetrachloroethene ND 1.0 0.29 1 2-Chlorotoluene ND 1.0 0.24 1 Toluene ND 1.0 0.35 1 4-Chlorotoluene ND 1.0 0.30 1 1,2,3-Trichlorobenzene ND 1.0 0.39 1 1,2-Dibromochloromethane ND 1.0 0.45 1 1,2,4-Trichloroethene ND 1.0 0.35 1 1,2-Dibromochloromethane ND 1.0 0.45 1 1,1,1-Trichloroethane ND 1.0 0.32 1 1,2-Dibromoethane ND 1.0 0.81 1 1,1,1-Trichloroethane ND 1.0 0.32 1 1,2-Dibromoethane ND 1.0 0.81 1 1,1,2-Trichloro-1,2,2-Trifluoroethane ND 1.0 0.54 1 1,2-Dichlorobenzene ND 1.0 0.42 1 1,1,2-Trichloroethane ND 1.0 0.54 1 1,2-Dichlorobenzene ND 1.0 0.38 1 Trichloroethene ND 1.0 0.30 1 1,3-Dichlorobenzene ND 1.0 0.38 1 Trichlorofluoromethane ND 1.0 0.36 1 1,4-Dichlorobenzene ND 1.0 0.30 1 1,2,3-Trichloropopane ND 1.0 0.36 1 1,1-Dichloroethane ND 1.0 0.53 1 1,2,3-Trimethylbenzene ND 1.0 0.26 1 1,1-Dichloroethane ND 1.0 0.53 1 1,3,5-Trimethylbenzene ND 1.0 0.26 1 1,1-Dichloroethane ND 1.0 0.31 1 Vinyl Chloride ND 0.50 0.33 1 1,1-Dichloroethene ND 1.0 0.35 1 Pym-Xylene ND 1.0 0.38 1 1,2-Dichloroethene ND 1.0 0.29 1 0.00 0.29 1 1,2-Dichloroethene ND 1.0 0.29 1 0.00 0.29 1 1,2-Dichloroethene ND 1.0 0.29 1 0.00 0.29 1 1,2-Dichloroethene ND 1.0 0.29 1 0.00 0.20 0.20 1 0.00 0.20 0.20 1 0.00 0.20 0.2	Chloroethane	ND	1.0	0.52	1		1,1,1,2-Tetrach	loroethane	ND	1.0	0.37	1	
2-Chlorotoluene ND 1.0 0.24 1 Toluene ND 1.0 0.35 1 4-Chlorotoluene ND 1.0 0.30 1 1,2,3-Trichlorobenzene ND 1.0 0.39 1 Dibromochloromethane ND 1.0 0.45 1 1,2,4-Trichlorobenzene ND 1.0 0.35 1 1,2-Dibromochloromethane ND 1.0 0.45 1 1,2,4-Trichlorobenzene ND 1.0 0.35 1 1,2-Dibromochlane ND 1.0 0.81 1 1,1,2-Trichloroch1,2,2-Trifluoroethane ND 1.0 0.32 1 1,2-Dibrlorobenzene ND 1.0 0.42 1 1,1,2-Trichloroethane ND 1.0 0.54 1 1,2-Dichlorobenzene ND 1.0 0.24 1 Trichloroethane ND 1.0 0.36 1 1,4-Dichlorobenzene ND 1.0 0.33 1 Trichloroethane ND 1.0 0.36 <td< td=""><td>Chloroform</td><td>ND</td><td>1.0</td><td>0.22</td><td>1</td><td></td><td>1,1,2,2-Tetrach</td><td>loroethane</td><td>ND</td><td>1.0</td><td>0.37</td><td>1</td><td></td></td<>	Chloroform	ND	1.0	0.22	1		1,1,2,2-Tetrach	loroethane	ND	1.0	0.37	1	
4-Chlorotoluene ND 1.0 0.30 1 1,2,3-Trichlorobenzene ND 1.0 0.39 1 1,2-Dibromo-3-Chloropropane ND 5.0 2.5 1 1,1,1-Trichloroethane ND 1.0 0.32 1 1,2-Dibromoethane ND 1.0 0.81 1 1,1,2-Trichloroethane ND 1.0 0.32 1 1,2-Dibromoethane ND 1.0 0.81 1 1,1,2-Trichloroethane ND 1.0 0.54 1 1,2-Dichloroethane ND 1.0 0.30 1 1,2,3-Trichloroethane ND 1.0 0.36 1 1,4-Dichloroethane ND 1.0 0.38 1 1 1,2,3-Trichloropropane ND 1.0 0.36 1 1,2,3-Trichloropropane ND 1.0 0.36 1 1,2,3-Trichloropropane ND 1.0 0.36 1 1,2,3-Trichloropropane ND 1.0 0.26 1 1,1-Dichloroethane ND 1.0 0.53 1 1,2,3-Trichloropropane ND 1.0 0.26 1 1,1-Dichloroethane ND 1.0 0.53 1 1,3,5-Trimethylbenzene ND 1.0 0.26 1 1,1-Dichloroethane ND 1.0 0.31 1 Vinyl Acetate ND 1.0 0.30 1 1,1-Dichloroethene ND 1.0 0.35 1 Vinyl Acetate ND 1.0 0.38 1 1,2-Dichloroethene ND 1.0 0.35 1 P/m-Xylene ND 1.0 0.29 1 1,2-Dichloroethene ND 1.0 0.29 1 0-Xylene ND 1.0 0.29 1 1,2-Dichloroethene ND 1.0 0.29 1 0-Xylene ND 1.0 0.29 1 1,2-Dichloroethene ND 1.0 0.29 1 0-Xylene ND 1.0 0.29 1 1,2-Dichloroethane ND 1.0 0.29 1 0-Xylene ND 1.0 0.29 1 1,2-Dichloropropane ND 1.0 0.29 1 0-Xylene ND 1.0 0.29 1 1,2-Dichloroethene ND 1.0 0.28 1 0-Xylene ND 1.0 0.29 1 1,2-Dichloroethane ND 1.0 0.28 1 0-Xylene ND 1.0 0.29 1 1,2-Dichloroethane ND 1.0 0.28 1 0-Xylene ND 1.0 0.29 1 1,2-Dichloroethane ND 1.0 0.28 1 0-Xylene ND 1.0 0.29 1 1,2-Dichloroethane ND 1.0 0.29 1 0-Xylene ND 1.0 0.29 1 1,2-Dichloroethane ND 1.0 0.29 1 0-Xylene ND 1.0 0.29 1 1,2-Dichloroethane ND 1.0 0.29 1 0-Xylene ND 1.0 0.29 1 1,2-Dichloroethane ND 1.0 0.29 1 0-Xylene ND 1.0 0.29 1 1,2-Dichloroethane ND 1.0 0.29 1 0-Xylene ND 1.0 0.29 1 1,2-Dichloroethane ND 1.0 0.29 1 0-Xylene ND 1.0 0.29 1 1,2-Dichloroethane ND 1.0 0.29 1 0-Xylene ND 1.0 0.29 1	Chloromethane	ND	10	1.8	1		Tetrachloroethe	ene	ND	1.0	0.29	1	
A-Chlorotoluene ND 1.0 0.30 1 1,2,3-Trichlorobenzene ND 1.0 0.39 1	2-Chlorotoluene	ND	1.0	0.24	1		Toluene		ND	1.0	0.35	1	
1,2-Dibromo-3-Chloropropane ND 5.0 2.5 1 1,1,1-Trichloroethane ND 1.0 0.32 1 1,2-Dibromoethane ND 1.0 0.81 1 1,1,2-Trichloroethane ND 10 0.54 1 Dibromomethane ND 1.0 0.42 1 1,1,2-Trichloroethane ND 1.0 0.54 1 1,2-Dichlorobenzene ND 1.0 0.24 1 Trichloroethene ND 1.0 0.30 1 1,3-Dichlorobenzene ND 1.0 0.38 1 Trichloroethene ND 10 0.36 1 1,4-Dichlorobenzene ND 1.0 0.30 1 1,2,3-Trichloropropane ND 5.0 2.3 1 Dichloroethane ND 1.0 0.27 1 1,2,4-Trimethylbenzene ND 1.0 0.26 1 1,1-Dichloroethane ND 1.0 0.53 1 1,3,5-Trimethylbenzene ND 1.0 0.32 1		ND	1.0	0.30	1		1,2,3-Trichlorob	benzene	ND	1.0	0.39	1	
1,2-Dibromoethane ND 1.0 0.81 1 1,1,2-Trichloro-1,2,2-Trifluoroethane ND 10 0.54 1 Dibromomethane ND 1.0 0.42 1 1,1,2-Trichloroethane ND 1.0 0.54 1 1,2-Dichlorobenzene ND 1.0 0.24 1 Trichloroethene ND 1.0 0.30 1 1,3-Dichlorobenzene ND 1.0 0.38 1 Trichlorofluoromethane ND 10 0.36 1 1,4-Dichlorobenzene ND 1.0 0.30 1 1,2,3-Trichloropropane ND 5.0 2.3 1 Dichlorodifluoromethane ND 1.0 0.27 1 1,2,4-Trimethylbenzene ND 1.0 0.26 1 1,1-Dichloroethane ND 1.0 0.53 1 1,3,5-Trimethylbenzene ND 1.0 0.19 1 1,2-Dichloroethane ND 1.0 0.31 1 Vinyl Acetate ND 1.0 0.30	Dibromochloromethane	ND	1.0	0.45	1		1,2,4-Trichlorob	benzene	ND	1.0	0.35	1	
Dibromomethane	1,2-Dibromo-3-Chloropropane	ND	5.0	2.5	1		1,1,1-Trichloroe	ethane	ND	1.0	0.32	1	
1,2-Dichlorobenzene	1,2-Dibromoethane	ND	1.0	0.81	1		1,1,2-Trichloro-	1,2,2-Trifluoroet	hane ND	10	0.54	1	
1,3-Dichlorobenzene ND 1.0 0.38 1 Trichlorofluoromethane ND 10 0.36 1 1,4-Dichlorobenzene ND 1.0 0.30 1 1,2,3-Trichloropropane ND 5.0 2.3 1 Dichlorodifluoromethane ND 1.0 0.27 1 1,2,4-Trimethylbenzene ND 1.0 0.26 1 1,1-Dichloroethane ND 1.0 0.53 1 1,3,5-Trimethylbenzene ND 1.0 0.19 1 1,2-Dichloroethane ND 0.50 0.22 1 Vinyl Acetate ND 1.0 0.19 1 1,1-Dichloroethene ND 1.0 0.31 1 Vinyl Acetate ND 1.0 0.32 1 1,1-Dichloroethene ND 1.0 0.31 1 Vinyl Chloride ND 0.50 0.33 1 c-1,2-Dichloroethene ND 1.0 0.29 1 o-Xylene ND 1.0 0.21 1	Dibromomethane	ND	1.0	0.42	1		1,1,2-Trichloroe	ethane	ND	1.0	0.54	1	
1,4-Dichloroberzene ND 1.0 0.30 1 1,2,3-Trichloropropane ND 5.0 2.3 1 Dichlorodifluoromethane ND 1.0 0.27 1 1,2,4-Trimethylbenzene ND 1.0 0.26 1 1,1-Dichloroethane ND 1.0 0.53 1 1,3,5-Trimethylbenzene ND 1.0 0.19 1 1,2-Dichloroethane ND 0.50 0.22 1 Vinyl Acetate ND 10 3.2 1 1,1-Dichloroethane ND 1.0 0.31 1 Vinyl Acetate ND 10 3.2 1 1,1-Dichloroethene ND 1.0 0.31 1 Vinyl Chloride ND 0.50 0.33 1 c-1,2-Dichloroethene ND 1.0 0.35 1 p/m-Xylene ND 1.0 0.38 1 t-1,2-Dichloroethene ND 1.0 0.29 1 o-Xylene ND 1.0 0.29 1	1,2-Dichlorobenzene	ND	1.0	0.24	1		Trichloroethene)	ND	1.0	0.30	1	
Dichlorodifluoromethane ND 1.0 0.26 1 1,1-Dichloroethane ND 1.0 0.53 1 1,3,5-Trimethylbenzene ND 1.0 0.19 1 1,2-Dichloroethane ND 0.50 0.22 1 Vinyl Acetate ND 10 3.2 1 1,1-Dichloroethane ND 1.0 0.31 1 Vinyl Chloride ND 0.50 0.33 1 c-1,2-Dichloroethane ND 1.0 0.35 1 p/m-Xylene ND 1.0 0.38 1 t-1,2-Dichloroethane ND 1.0 0.29 1 o-Xylene ND 1.0 0.21 1 1,2-Dichloropropane ND 1.0 0.28 1 Methyl-t-Butyl Ether (MTBE) ND 1.0 0.29 1 Surrogates: REC (%) Control Limits Qual Surrogates: REC (%) Control Limits Qual	1,3-Dichlorobenzene	ND	1.0	0.38			Trichlorofluoron	nethane	ND	10	0.36	1	
1,1-Dichloroethane ND 1.0 0.53 1 1,3,5-Trimethylbenzene ND 1.0 0.19 1 1,2-Dichloroethane ND 0.50 0.22 1 Vinyl Acetate ND 10 3.2 1 1,1-Dichloroethene ND 1.0 0.31 1 Vinyl Chloride ND 0.50 0.33 1 c-1,2-Dichloroethene ND 1.0 0.35 1 p/m-Xylene ND 1.0 0.38 1 t-1,2-Dichloroethene ND 1.0 0.29 1 o-Xylene ND 1.0 0.21 1 1,2-Dichloropropane ND 1.0 0.28 1 Methyl-t-Butyl Ether (MTBE) ND 1.0 0.29 1 Surrogates: REC (%) Control Limits Qual Surrogates: REC (%) Control Limits Qual	1,4-Dichlorobenzene	ND	1.0	0.30	1		1,2,3-Trichlorop	oropane	ND	5.0	2.3	1	
1,2-Dichloroethane ND 0.50 0.22 1 Vinyl Acetate ND 10 3.2 1 1,1-Dichloroethene ND 1.0 0.31 1 Vinyl Chloride ND 0.50 0.33 1 c-1,2-Dichloroethene ND 1.0 0.35 1 p/m-Xylene ND 1.0 0.38 1 t-1,2-Dichloroethene ND 1.0 0.29 1 o-Xylene ND 1.0 0.21 1 1,2-Dichloropropane ND 1.0 0.28 1 Methyl-t-Butyl Ether (MTBE) ND 1.0 0.29 1 Surrogates: REC (%) Control Limits Qual Surrogates: REC (%) Control Limits Qual	Dichlorodifluoromethane	ND							ND	1.0	0.26	1	
1,1-Dichloroethene ND 1.0 0.31 1 Vinyl Chloride ND 0.50 0.33 1 c-1,2-Dichloroethene ND 1.0 0.35 1 p/m-Xylene ND 1.0 0.38 1 t-1,2-Dichloroethene ND 1.0 0.29 1 o-Xylene ND 1.0 0.21 1 1,2-Dichloropropane ND 1.0 0.28 1 Methyl-t-Butyl Ether (MTBE) ND 1.0 0.29 1 Surrogates: REC (%) Control Limits Qual Surrogates: REC (%) Control Limits Qual Dibromofluoromethane 116 74-140 1,2-Dichloroethane-d4 114 74-146 74-146	1,1-Dichloroethane	ND	1.0	-			1,3,5-Trimethyl	benzene	ND	- 1.0	0.19	1	
c-1,2-Dichloroethene ND 1.0 0.35 1 p/m-Xylene ND 1.0 0.38 1 t-1,2-Dichloroethene ND 1.0 0.29 1 o-Xylene ND 1.0 0.21 1 1,2-Dichloropropane ND 1.0 0.28 1 Methyl-t-Butyl Ether (MTBE) ND 1.0 0.29 1 Surrogates: REC (%) Control Limits Qual Surrogates: REC (%) Control Limits Qual Dibromofluoromethane 116 74-140 1,2-Dichloroethane-d4 114 74-146	1,2-Dichloroethane	ND	0.50				Vinyl Acetate		ND	10		1	
t-1,2-Dichloroethene ND 1.0 0.29 1 o-Xylene ND 1.0 0.21 1 1,2-Dichloropropane ND 1.0 0.28 1 Methyl-t-Butyl Ether (MTBE) ND 1.0 0.29 1 Surrogates: REC (%) Control Limits Qual Surrogates: REC (%) Control Limits Qual Dibromofluoromethane 116 74-140 1,2-Dichloroethane-d4 114 74-146		ND.	1.0		-		•		ND	-		•	
1,2-Dichloropropane ND 1.0 0.28 1 Methyl-t-Butyl Ether (MTBE) ND 1.0 0.29 1 Surrogates: REC (%) Control Limits Qual Surrogates: REC (%) Control Limits Qual Dibromofluoromethane 116 74-140 1,2-Dichloroethane-d4 114 74-146	c-1,2-Dichloroethene			_	1		p/m-Xylene		ND	1.0			
Surrogates: REC (%) Control Limits Qual Surrogates: REC (%) Control Limits Qual Dibromofluoromethane 116 74-140 1,2-Dichloroethane-d4 114 74-146	t-1,2-Dichloroethene	ND			1		•					•	
Dibromofluoromethane 116 74-140 1,2-Dichloroethane-d4 114 74-146					•	_		Ether (MTBE)				•	
	Surrogates:	<u>REC_(%)</u>		<u>imits</u>		<u>Qual</u>	Surrogates:		<u>REC (%</u>		<u>Limits</u>	Q	<u>ual</u>
Toluene-d8 99 88-112 1,4-Bromofluorobenzene 80 74-110	Dibromofluoromethane	116	74-140				1,2-Dichloroeth	ane-d4	114	74-146			
	Toluene-d8	99	88-112				1,4-Bromofluore	obenzene	80	74-110			







Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: Units:

Date

Date

Date

06-06-1183 EPA 5030B EPA 8260B ug/L

06/19/06

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

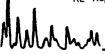
Lab Sample

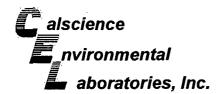
Page 2 of 8

Client Sample Number			Numi	ber	Collected	Matrix	Prepared	Analyzed	QC Bate	ch ID	
MW-6		1	06-06	1183-2	06/19/06	COLORS OF THE PARTY OF THE PART	06/22/06	06/23/06	060622	L02	
Comment(s): -Results were e	valuated to the	e MDL, co	ncentratio	ons >= to the I	MDL but < RL, if fo	ound, are qualifie	ed with a "J" f	lag.			
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>MDL</u>	DF Qual	<u>Parameter</u>		<u>Result</u>	<u>RL</u>	<u>MDL</u>	DF !	Qual
Acetone	ND	10	6.1	1	1,3-Dichloroprop	pane	ND	1.0	0.30	1	
Benzene	ND	0.50	0.26	1	2.2-Dichloropror	pane	ND	1.0	0.40	1	
Bromobenzene	ND	1.0	0.47	1	1,1-Dichloropro		ND	1.0	0.21	1	
Bromochloromethane	ND	1.0	0.68	1	c-1,3-Dichloropi	ropene	ND	0.50	0.45	1	
Bromodichloromethane	ND	1.0	0.27	1	t-1,3-Dichloropre	opene	ND	0.50	0.31	1	
Bromoform	ND	1.0	0.62	1	Ethylbenzene	•	ND	1.0	0.17	1	
Bromomethane	ND	10	2.9	1	2-Hexanone		ND '	10	1.9	1	
2-Butanone	ND	10	4.2	1	Isopropylbenzer	ne	ND	1.0	0.24	1	
n-Butylbenzene	ND	1.0	0.29	1	p-Isopropyltolue	ene	ND	1.0	0.21	1	
sec-Butylbenzene	ND	1.0	0.21	1	Methylene Chlor	ride	3.9	10.0	2.6	1	J,B
tert-Butylbenzene	ND	1.0	0.17	1	4-Methyl-2-Pent	tanone	ND	10	2.4	1	_
Carbon Disulfide	ND	10	1.0	1	Naphthalene		ND	10	0.95	1	4
Carbon Tetrachloride	3.1	0.5	0.42	1	n-Propylbenzen	е	ND	1.0	0.30	1	•
Chlorobenzene	ND	1.0	0.36	1	Styrene		ND	1.0	0.29	1	
Chloroethane	ND	1.0	0.52	1	1,1,1,2-Tetrachi	loroethane	ND	1.0	0.37	1	
Chloroform	2.4	1.0	0.22	1	1,1,2,2-Tetrachl	loroethane	ND	1.0	0.37	1	
Chloromethane	ND	10	1.8	1	Tetrachloroethe	ne	150	1	0.29	1	
2-Chlorotoluene	ND	1.0	0.24	1	Toluene		ND	1.0	0.35	1	
4-Chlorotoluene	ND	1.0	0.30	1	1,2,3-Trichlorob	enzene	ND	1.0	0.39	1	
Dibromochloromethane	ND	1.0	0.45	1	1,2,4-Trichlorob	enzene	ND	1.0	0.35	1	
1,2-Dibromo-3-Chloropropane	ND	5.0	2.5	1	1,1,1-Trichloroe	ethane	ND	1.0	0.32	1	
1,2-Dibromoethane	ND	1.0	0.81	1∙	1,1,2-Trichloro-	1,2,2-Trifluoroet	hane 1.1	10.0	0.54	1	J
Dibromomethane	ND	1.0	0.42	1	1,1,2-Trichloroe	thane	ND	1.0	0.54	1	
1,2-Dichlorobenzene	ND	1.0	0.24	1	Trichloroethene		75	1	0.30	1	
1,3-Dichlorobenzene	ND	1.0	0.38	1	Trichlorofluorom	nethane	ND	10	0.36	1	
1,4-Dichlorobenzene	ND	1.0	0.30	.1	1,2,3-Trichlorop	ropane	ND	5.0	2.3	1	
Dichlorodifluoromethane	ND	1.0	0.27	1	1,2,4-Trimethylb	benzene	ND	1.0	0.26	1	
1,1-Dichloroethane	ND	1.0	0.53	1	1,3,5-Trimethylb	benzene	ND	1.0	0.19	1	
1,2-Dichloroethane	ND	0.50	0.22	1	Vinyl Acetate		ND	10	3.2	1	
1,1-Dichloroethene	3.4	1.0	0.31	1	Vinyl Chloride		ND	0.50	0.33	1	
c-1,2-Dichloroethene	ND	1.0	0.35	1	p/m-Xylene		ND	1.0	0.38	1	
t-1,2-Dichloroethene	ND	1.0	0.29	1	o-Xylene		ND	1.0	0.21	1	
1,2-Dichloropropane	ND	1.0	0.28	1	Methyl-t-Butyl E	ther (MTBE)	ND	1.0	0.29	1	
Surrogates:	<u>REC (%)</u>	Control	<u>Limits</u>	<u>Qual</u>	Surrogates:		<u>REC (%</u>) <u>Control</u>	<u>Limits</u>	:	Qual
Dibromofluoromethane	116	74-140			1,2-Dichloroetha	ane-d4	116	74-146			
Toluene-d8	100	88-112			1,4-Bromofiuoro	obenzene	79	74-110			

RL - Reporting Limit

DF - Dilution Factor







Tetra Tech, Inc.

Carbon Tetrachloride

Chlorobenzene Chloroethane

Chloromethane

2-Chlorotoluene

Chloroform

Toluene-d8

3475 East Foothill Blvd., Suite 300

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

0.5

1.0

1.0

1.0

1.0

10

2.4

2.0

ND

ND

ND

ND

100

0.42

0.36

0.52

0.22

1.8

0.24

Pasadena, CA 91107-6024

Date Received:

Work Order No:

Preparation:

Method:

n-Propylbenzene

Tetrachloroethene

1,1,1,2-Tetrachloroethane

1,1,2,2-Tetrachloroethane

Styrene

Toluene

Units:

06-06-1183 EPA 5030B

06/19/06

EPA 8260B

A 02000

ug/L Page 3 of 8

ND

ND

ND

ND

150

ND

78

1.0

1.0

1.0

1.0

1.0

74-110

1

0.30

0.29

0.37

0.37

0.29

0.35

Client Sample Number			Lab Sample Number			Date Collected Matrix	Date Prepared	Date Analyzed	QC Batch ID	
MW-5			06-06	1183-3		06/19/06 Aqueous	% 06/ 22/ 06	06/23/06	060622	L02
		ne MDL, co	oncentratio	ons >= 1	to the N	MDL but < RL, if found, are qu	ualified with a "J"	flag.		
<u>Parameter</u>	Result	RL	<u>MDL</u>	<u>DF</u>	Qual	<u>Parameter</u>	Resul	t <u>RL</u>	<u>MDL</u>	DF Qual
Acetone	6.3	10.0	6.1	1	J,B	1,3-Dichloropropane	, ND	1.0	0.30	1
Benzene	ND	0.50	0.26	1		2,2-Dichloropropane	ND	1.0	0.40	1
Bromobenzene	ND	1.0	0.47	1		1,1-Dichloropropene	ND	1.0	0.21	1
Bromochloromethane	ND	1.0	0.68	1		c-1,3-Dichloropropene	ND	0.50	0.45	1
Bromodichloromethane	0.29	1.0	0.27	1	J	t-1,3-Dichloropropene	ND	0.50	0.31	1
Bromoform	ND	1.0	0.62	1		Ethylbenzene	ND	1.0	0.17	1
Bromomethane	ND	10	2.9	1		2-Hexanone	ND	10	1.9	1
2-Butanone	ND -	10	4.2	1		Isopropylbenzene	ND	1.0	0.24	1
n-Butylbenzene	ND	1.0	0.29	1		p-Isopropyttoluene	ND	1.0	0.21	1
sec-Butylbenzene	ND	1.0	0.21	1		Methylene Chloride	4.6	10.0	2.6	1 J,B
tert-Butylbenzene	ND	1.0	0.17	1		4-Methyl-2-Pentanone	ND	10	2.4	լ1
Carbon Disulfide	ND	10	1.0	1		Naphthalene	ND	10	0.95	1

1

1

1

1

1

4-Chlorotoluene ND 1.0 0.30 1 1,2,3-Trichlorobenzene ND 1.0 0.39 1 1 0.45 1,2,4-Trichlorobenzene 0.35 1 ND 1.0 Dibromochloromethane ND 1.0 ND 5.0 2.5 1.1.1-Trichloroethane ND 1.0 0.32 1 1,2-Dibromo-3-Chloropropane 0.81 1,1,2-Trichloro-1,2,2-Trifluoroethane 10.0 0.54 1,2-Dibromoethane ND 1.0 1.6 0.54 1 Dibromomethane ND 1.0 0.42 1,1,2-Trichloroethane ND 1.0 Trichloroethene 0.30 1 ND 1.0 0.24 84 1,2-Dichlorobenzene 1 10 0.36 ND 1.0 0.38 Trichlorofluoromethane ND 1.3-Dichlorobenzene 1 1,4-Dichlorobenzene ND 1.0 0.30 1,2,3-Trichloropropane ND 5.0 2.3 1 1 Dichlorodifluoromethane ND 1.0 0.27 1,2,4-Trimethylbenzene ND 1.0 0.26 1 ND 1.0 0.53 1 1,3,5-Trimethylbenzene ND 1.0 0.19 1 1,1-Dichloroethane 0.22 0.29 0.50 1 Vinyl Acetate ND 10 3.2 1 1,2-Dichloroethane Vinvl Chloride 1.0 0.31 ND 0.50 0.33 1,1-Dichloroethene 2.6 c-1,2-Dichloroethene ND 1.0 0.35 1 p/m-Xylene ND 1.0 0.38 1 1 0.29 o-Xylene 1.0 1.0 ND 0.21 t-1,2-Dichloroethene ND ND 1.0 0.28 Methyl-t-Butyl Ether (MTBE) ND 1.0 0.29 1,2-Dichloropropane **REC (%)** Control Limits Surrogates: **REC (%)** Control Limits Qual Surrogates: 74-140 1,2-Dichloroethane-d4 124 74-146 Dibromofluoromethane 126

RL - Reporting Limit .

DF - Dilution Factor

88-112

Qual - Qualifiers

1,4-Bromofluorobenzene





Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: 06/19/06 06-06-1183 EPA 5030B EPA 8260B

Units:

ug/L

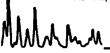
Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

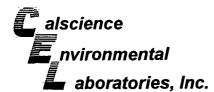
Page 4 of 8

Client Sample Number			Lab Sa Numb		Date Collected Matrix	Date Prepared	Date Analyzed	QC Bate	ch ID
MW4			06-06	1183-4		06/22/06	.,06/23/06	060622	02
Comment(s): -Results were e	evaluated to the	e MDL, co	ncentratio	ons >= to the N	ADL but < RL, if found, are qua	lified with a "J"	flag.		
<u>Parameter</u>	Result	<u>RL</u>	<u>MDL</u>	DF Qual	<u>Parameter</u>	Resu	it <u>RL</u>	MDL.	DF Qual
Acetone	ND	10	6.1	1	1,3-Dichloropropane	ND	1.0	0.30	1
Benzene	ND	0.50	0.26	1	2,2-Dichloropropane	ND	1.0	0.40	1
Bromobenzene	ND	1.0	0.47	1	1,1-Dichloropropene	ND	1.0	0.21	1
Bromochloromethane	ND	1.0	0.68	1	c-1,3-Dichloropropene	ND	0.50	0.45	1
Bromodichloromethane	ND	1.0	0.27	1	t-1,3-Dichloropropene	ND	0.50	0.31	1
Bromoform	ND	1.0	0.62	1	Ethylbenzene	ND	1.0	0.17	1
Bromomethane	ND	10	2.9	1	2-Hexanone	ND	10	1.9	1
2-Butanone	ND	10	4.2	1	Isopropylbenzene	ND	1.0	0.24	1
n-Butylbenzene	ND	1.0	0.29	1	p-isopropyttoluene	ND	1.0	0.21	1
sec-Butylbenzene	ND	1.0	0.21	1	Methylene Chloride	4.5	10.0	2.6	1 J,B
tert-Butylbenzene	ND	1.0	0.17	1	4-Methyl-2-Pentanone	ND	10	2.4	1
Carbon Disulfide	ND	10	1.0	1	Naphthalene	ND	10	0.95	1
Carbon Tetrachloride	0.81	0.50	0.42	1	n-Propylbenzene	ND	1.0	0.30	1
Chlorobenzene	ND	1.0	0.36	1	Styrene	ND	1.0	0.29	1
Chloroethane	ND	1.0	0.52	1	1,1,1,2-Tetrachloroethane	ND	1.0	0.37	1
Chloroform	1.3	1.0	0.22	1	1,1,2,2-Tetrachloroethane	ND	1.0	0.37	1
Chloromethane	ND	10	1.8	1	Tetrachloroethene	120	1	0.29	1
2-Chlorotoluene	ND	1.0	0.24	1	Toluene	ND	1.0	0.35	1
4-Chlorotoluene	ND	1.0	0.30	1	1,2,3-Trichlorobenzene	ND	1.0	0.39	1
Dibromochloromethane	ND	1.0	0.45	1	1,2,4-Trichlorobenzene	ND	1.0	0.35	1
1.2-Dibromo-3-Chloropropane	ND	5.0	2.5	1	1,1,1-Trichloroethane	ND	1.0	0.32	1
1.2-Dibromoethane	ND	1.0	0.81	1	1,1,2-Trichloro-1,2,2-Trifluoro	ethane 1.1	10.0	0.54	1 J
Dibromomethane	ND	1.0	0.42	1	1,1,2-Trichloroethane	ND	1.0	0.54	1
1,2-Dichlorobenzene	ND	1.0	0.24	1	Trichloroethene	48	1	0.30	1
1,3-Dichlorobenzene	ND	1.0	0.38	1	Trichlorofluoromethane	ND	10	0.36	1
1,4-Dichlorobenzene	ND	1.0	0.30	1	1,2,3-Trichloropropane	ND	5.0	2.3	1
Dichlorodifluoromethane	ND	1.0	0.27	1	1,2,4-Trimethylbenzene	ND	1.0	0.26	1
1,1-Dichloroethane	ND	1.0	0.53	1	1,3,5-Trimethylbenzene	ND	1.0	0.19	1
1,2-Dichloroethane	ND	0.50	0.22	1	Vinyl Acetate	ND	10	3.2	1
1,1-Dichloroethene	0.64	1.0	0.31	1 J	Vinyl Chloride	ND	0.50	0.33	1
c-1,2-Dichloroethene	ND	1.0	0.35	1	p/m-Xylene	ND	1.0	0.38	1
t-1,2-Dichloroethene	ND	1.0	0.29	1	o-Xylene	ND	1.0	0.21	1
1,2-Dichloropropane	ND	1.0	0.28	1	Methyl-t-Butyl Ether (MTBE)	ND	1.0	0.29	1
Surrogates:	REC (%)	Control	<u>Limits</u>	<u>Qual</u>	Surrogates:	REC (%) Contro	l Limits	<u>Qual</u>
Dibromofluoromethane	125	74-140			1,2-Dichloroethane-d4	126	74-146		
Toluene-d8	99	88-112			1,4-Bromofluorobenzene	78	74-110		

RL - Reporting Limit

DF - Dilution Factor







Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

06/19/06

Work Order No:

06-06-1183

Preparation:

EPA 5030B

Method:

EPA 8260B

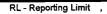
Units:

ug/L

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

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Client Sample Number				Lab Sample Number		Matrix	Date Prepared	Date Analyzed	QC Bat	ch ID	
MW4D	7157		06-06	1183-5	06/19/06	Aqueous	06/23/06	06/24/06	060623	L02:	ŝΫ÷
Comment(s): -Results were e	evaluated to the	MDL, co	ncentrati	ons >= to the	MDL but < RL, if	found, are qualif	ied with a "J" t	Nag.			
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>MDL</u>	DF Qua	<u>Parameter</u>		Result	<u>RL</u>	MDL	<u>D</u> F	Qual
Acetone	9.5	10.0	6.1	1.	1,3-Dichloropi	ropane	ND	1.0	0.30	1	
Benzene	ND	0.50	0.26	1	2,2-Dichloropi		ND	1.0	0.40	1	
Bromobenzene	ND	1.0	0.47	1	1,1-Dichloropa	ropene	ND	1.0	0.21	1	
Bromochloromethane	ND	1.0	0.68	1	c-1,3-Dichloro	propene	ND	0.50	0.45	1	
Bromodichloromethane	ND	1.0	0.27	1	t-1,3-Dichloro	propene	ND	0.50	0.31	1	
Bromoform	ND	1.0	0.62	1	Ethylbenzene		ND	1.0	0.17	1	
Bromomethane	ND	10	2.9	1	2-Hexanone		ND	10	1.9	1	
2-Butanone	ND	10	4.2	1	Isopropylbenz	ene	ND	1.0	0.24	1	
n-Butylbenzene	ND	1.0	0.29	1	p-Isopropyltoli	uene	ND	1.0	0.21	1	
sec-Butylbenzene	ND	1.0	0.21	1	Methylene Ch	loride	2.9	10.0	2.6	1	J,B
tert-Butylbenzene	ND	1.0	0.17	1	4-Methyl-2-Pe	entanone	ND	10	2.4	1	
Carbon Disulfide	ND	10	1.0	1	Naphthalene		ND	10	0.95	1	
Carbon Tetrachloride	0.88	0.50	0.42	1	n-Propylbenze	ene	ND	1.0	0.30	1	
Chlorobenzene	ND	1.0	0.36	1	Styrene		ND	1.0	0.29	1	
Chloroethane	ND	1.0	0.52	1	1,1,1,2-Tetrac	chloroethane	ND	1.0	0.37	1	
Chloroform	1.2	1.0	0.22	1	1,1,2,2-Tetrac	chloroethane	ND	1.0	0.37	1	
Chloromethane	ND	10	1.8	1	Tetrachloroeth	nene	120	1	0.29	1	
2-Chlorotoluene	ND	1.0	0.24	1	Toluene		ND	1.0	0.35	1	
4-Chiorotoluene	ND	1.0	0.30	1	1,2,3-Trichlore	obenzene	ND	1.0	0.39	1	
Dibromochloromethane	ND	1.0	0.45	1	1,2,4-Trichlore	obenzene	ND	1.0	0.35	1	
1,2-Dibromo-3-Chloropropane	ND	5.0	2.5	1	1,1,1-Trichlore	oethane	ND	1.0	0.32	1	
1,2-Dibromoethane	ND	1.0	0.81	1	1,1,2-Trichlore	o-1,2,2-Trifluoroe	thane 1.0	10.0	0.54	1	J
Dibromomethane	ND	1.0	0.42	1	1,1,2-Trichlore	oethane	ND	1.0	0.54	1	
1,2-Dichlorobenzene	ND	1.0	0.24	1	Trichloroether	ne	48	1	0.30	1	
1,3-Dichlorobenzene	ND	1.0	0.38	1	Trichlorofluoro	omethane	ND	10	0.36	1	
1,4-Dichlorobenzene	ND	1.0	0.30	1	1,2,3-Trichlore	opropane	ND	5.0	2.3	1	
Dichlorodifluoromethane	ND	1.0	0.27	1	1,2,4-Trimethy	ylbenzene	, ND	1.0	0.26	1	
1,1-Dichloroethane	ND	1.0	0.53	1	1,3,5-Trimethy	ylbenzene	ND	1.0	0.19	1	
1,2-Dichloroethane	ND	0.50	0.22	1	Vinyl Acetate		ND	10	3.2	1	
1,1-Dichloroethene	0.82	1.0	0.31		Vinyl Chloride		ND	0.50	0.33	1	
c-1,2-Dichloroethene	ND	1.0	0.35	1	p/m-Xylene		ND	1.0	0.38	1	
t-1,2-Dichloroethene	ND	1.0	0.29	1	o-Xylene		ND	1.0	0.21	1	
1,2-Dichloropropane	ND	1.0	0.28	1		Ether (MTBE)	ND	1.0	0.29	1	_
Surrogates:	REC (%)	Control	<u>Limits</u>	Qua	·		REC (9		Limits		Qual
Dibromofluoromethane	123	74-140			1,2-Dichloroet	thane-d4	120	74-146			
Toluene-d8	98	88-112			1,4-Bromofluo	orobenzene	79	74-110			







Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Date Received: Work Order No: Preparation:

Date

Date

06/19/06 06-06-1183 **EPA 5030B**

Method: Units:

EPA 8260B

Date

ug/L Page 6 of 8

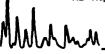
Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

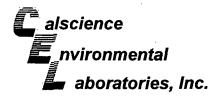
Lab Sample

OC Batch ID

Client Sample Number			Numl	Number Collected Matrix		Prepared	Analyzed	QC Bate	ch ID			
LFB061906		Í.	06-06-	1183-6		06/19/06; ** Aqueous -> 0		06/22/06	06/23/06	0606221	L02	
Comment(s): -Results were e	evaluated to the	e MDL, co	DL, concentrations >= to the MDL but < RL, if found, are quali		fied with a "J" f	lag.						
<u>Parameter</u>	Result	<u>RL</u>	<u>MDL</u>	<u>DF</u>	Qual	<u>Parameter</u>		Result	<u>RL</u>	MDL	<u>DF</u>	Qual
Acetone	14	10	6.1	1	В	1,3-Dichloropro	opane	ND	1.0	0.30	1	
Benzene	ND	0.50	0.26	1		2,2-Dichloropro	pane	ND	1.0	0.40	1	
Bromobenzene	ND	1.0	0.47	1		1,1-Dichloropro	opene	ND	1.0	0.21	1	
Bromochloromethane	ND	1.0	0.68	1		c-1,3-Dichloro	propene	ND	0.50	0.45	1	
Bromodichloromethane	ND	1.0	0.27	1		t-1,3-Dichlorop	ropene	ND	0.50	0.31	1	
Bromoform	ND	1.0	0.62	1		Ethylbenzene		ND	1.0	0.17	1	
Bromomethane	ND	10	2.9	1		2-Hexanone		ИD	10	1.9	1	
2-Butanone	ND	10	4.2	1		Isopropylbenze	ene	ND	1.0	0.24	1	
n-Butylbenzene	ND	1.0	0.29	1		p-Isopropyltolu	ene	ND	1.0	0.21	1	
sec-Butylbenzene	ND	1.0	0.21	1		Methylene Chle	4.5	10.0	2.6	1	J,B	
tert-Butylbenzene	ND	1.0	0.17	1		4-Methyl-2-Per	ntanone	ND	10	2.4	1	
Carbon Disulfide	ND	10	1.0	1		Naphthalene		ND	10	0.95	1	4
Carbon Tetrachloride	ND	0.50	0.42	1		n-Propylbenze	ne	ND	1.0	0.30	1	•
Chlorobenzene	ND	1.0	0.36	1		Styrene		ND	1.0	0.29	1	
Chloroethane	ND	1.0	0.52	1		1,1,1,2-Tetracl	hloroethane	ND	1.0	0.37	1	
Chloroform	ND	1.0	0.22	1		1,1,2,2-Tetracl	hloroethane	ND	1.0	0.37	1	
Chloromethane	ND	10	1.8	1		Tetrachloroeth		ND	1.0	0.29	1	
2-Chiorotoluene	ND	1.0	0.24	1		Toluene		ND	1.0	0.35	1	
4-Chlorotoluene	ND	1.0	0.30	1		1,2,3-Trichloro	benzene	ND	1.0	0.39	1	
Dibromochloromethane	ND	1.0	0.45	1		1,2,4-Trichloro	benzene	ND	1.0	0.35	1	
1,2-Dibromo-3-Chloropropane	ND	5.0	2.5	1		1,1,1-Trichloro	ethane	ND	1.0	0.32	1	
1,2-Dibromoethane	ND	1.0	0.81	1		1,1,2-Trichloro	-1,2,2-Trifluoroe	ethane ND	· 10	0.54	1	
Dibromomethane	ND	1.0	0.42	1		1,1,2-Trichloro	ethane	ND	1.0	0.54	1	
1,2-Dichlorobenzene	ND	1.0	0.24	1		Trichloroethen	е	ND	1.0	0.30	-1	
1,3-Dichlorobenzene	NĐ	1.0	0.38	1		Trichlorofluoro	methane	ND	10	0.36	1	
1,4-Dichlorobenzene	ND	1.0	0.30	1		1,2,3-Trichloro	propane	ND	5.0	2.3	1	
Dichlorodifluoromethane	ND	1.0	0.27	1		1,2,4-Trimethy	lbenzene	ND	1.0	0.26	1	
1,1-Dichloroethane	ND ·	1.0	0.53	1		1,3,5-Trimethy	lbenzene	ND	1.0	0.19	1	
1,2-Dichloroethane	ND	0.50	0.22	1		Vinyl Acetate		· ND	10	3.2	1	
1,1-Dichloroethene	ND	1.0	0.31	1		Vinyl Chloride		ND	0.50	0.33	1	
c-1,2-Dichloroethene	ND	1.0	0.35	1		p/m-Xylene		ND	1.0	0.38	1	
t-1,2-Dichloroethene	ND	1.0	0.29	1		o-Xylene		ND	1.0	0.21	1	
1,2-Dichloropropane	ND	1.0	0.28	1		Methyl-t-Butyl	Ether (MTBE)	ND	1.0	0.29	1	
Surrogates:	REC (%)	Control	<u>Limits</u>		Qual	Surrogates:		<u>REC (%</u>	<u>Control</u>	<u>Limits</u>		<u>Qual</u>
Dibromofluoromethane	123	74-140				1,2-Dichloroeth	nane-d4	120	74-146			
Toluene-d8	98	88-112				1,4-Bromofluo	robenzene	81	74-110			

DF - Dilution Factor







Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

06/19/06

Work Order No:

Date

Date

06-06-1183

Preparation:

EPA 5030B

Method:

Date

EPA 8260B

Units:

ug/L

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

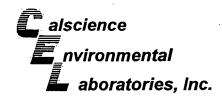
Lab Sample

Page 7 of 8

Client Sample Number			Number		Collected	Matrix	Prepared	Analyzed	QC Bate	ch ID
Method Blank	4 (1)	37.94	099-10	0-006-18,334	N/A	Aqueous	06/22/06	06/23/06	060622	102
	A SECTION OF THE PROPERTY.		Mary Secretary Control				(1 the street and the Carlot	regularistas (200)	Walter State Control	V///22/88#419#
Comment(s): -Results were ev	aluated to the	e MDL, cor	ncentratio	ons >= to the N	ADL but < RL, if	found, are qualif	ied with a "J" fl	ag.		
<u>Parameter</u>	Result	<u>RL</u>	<u>MDL</u>	DF Qual	<u>Parameter</u>		Result	<u>RL</u>	MDL	DF Qual
Acetone	9.8	10.0	6.1	1 J	1,3-Dichloropr	opane	ND	1.0	0.30	1
Benzene	ND	0.50	0.26	1	2,2-Dichloropr		ND	1.0	0.40	1
Bromobenzene	ND	1.0	0.47	1	1,1-Dichloropr	opene	ND	1.0	0.21	1
Bromochloromethane	ND	1.0	0.68	1	c-1,3-Dichloro		ND	0.50	0.45	1
Bromodichloromethane	ND	1.0	0.27	1	t-1,3-Dichlorop	propene	ND	0.50	0.31	1
Bromoform	ND	1.0	0.62	1 '	Ethylbenzene	•	ND	1.0	0.17	1
Bromomethane	ND	10	2.9	1	2-Hexanone		ND	10	1.9	1
2-Butanone	ND	10	4.2	1	Isopropylbenze	ene	ND	1.0	0.24	1
n-Butylbenzene	ND	1.0	0.29	1	p-Isopropyttolu	uene	ND	1.0	0.21	1
sec-Butylbenzene	ND	1.0	0.21	1	Methylene Chl	loride	4.1	10.0	2.6	1 J
tert-Butylbenzene	ND	1.0	0.17	1	4-Methyl-2-Pe	ntanone	ND	10	2.4	1
Carbon Disulfide	ND	10	1.0	1	Naphthalene		ND	10	0.95	1
Carbon Tetrachloride	ND	0.50	0.42	1	n-Propylbenze	ene	ND	1.0	0.30	1
Chlorobenzene	ND	1.0	0.36	1	Styrene		ND	1.0	0.29	1
Chloroethane	ND	1.0	0.52	1	1,1,1,2-Tetrac	hloroethane	ND	1.0	0.37	1
Chloroform	ND	1.0	0.22	1	1,1,2,2-Tetrac	hloroethane	ND	1.0	0.37	1
Chloromethane	ND	10	1.8	1	Tetrachloroeth	iene	ND	1.0	0.29	1
2-Chlorotoluene	ND	1.0	0.24	1	Toluene	•	ND	1.0	0.35	1
4-Chlorotoluene	ND	1.0	0.30	1	1,2,3-Trichlord	benzene	ND	1.0	0.39	1
Dibromochloromethane	ND	1.0	0.45	1	1,2,4-Trichlord	benzene	ND	1.0	0.35	1
1,2-Dibromo-3-Chloropropane	ND	5.0	2.5	1	1,1,1-Trichlord	ethane	ND	1.0	0.32	1
1,2-Dibromoethane	ND	1.0	0.81	1	1,1,2-Trichlord	-1,2,2-Trifluoroe	thane ND	10	0.54	1
Dibromomethane	ND	1.0	0.42	1	1,1,2-Trichlord	ethane	ND	1.0	0.54	1
1,2-Dichlorobenzene	ND	1.0	0.24	1	Trichloroethen	e	ND	1.0	0.30	1
1,3-Dichlorobenzene	ND	1.0	0.38	1	Trichlorofluoro		ND	10	0.36	1
1,4-Dichlorobenzene	ND	1.0	0.30	1 .	1,2,3-Trichlord		ND	5.0	2.3	1
Dichlorodifluoromethane	ND	1.0	0.27	1	1,2,4-Trimethy		ND	1.0	0.26	1
1,1-Dichloroethane	ND	1.0	0.53	1	1,3,5-Trimethy	/ibenzene	ND	1.0	0.19	1
1,2-Dichloroethane	ND	0.50	0.22	1	Vinyl Acetate		ND	10	3.2	1
1,1-Dichloroethene	ND	1.0	0.31	1	Vinyl Chloride		ND	0.50	0.33	1
c-1,2-Dichloroethene	ND	1.0	0.35	1	p/m-Xylene		ND	1.0	0.38	1
t-1,2-Dichloroethene	ND	1.0	0.29	1	o-Xylene	44-5-	ND	1.0	0.21	1
1,2-Dichloropropane	ND	1.0	0.28	1 Oual	Methyl-t-Butyl	Lther (MTBE)	ND DEC (9/	1.0	0.29	1
Surrogates:	REC (%)	Control L	<u>.imits</u>	<u>Qual</u>	Surrogates:		<u>REC (%</u>	_	LIMITS	<u>Qual</u>
Dibromofluoromethane	112	74-140			1,2-Dichloroet		110	74-146		
Toluene-d8	97	88-112			1,4-Bromofluo	robenzene	81	74-110		



DF - Dilution Factor





Date

Date

Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

Work Order No:

Preparation: Method:

Units:

Date

06/19/06

06-06-1183 **EPA 5030B**

EPA 8260B

ug/L

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

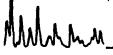
Lab Sample

Page 8 of 8

Client Sample Number			Lab Sample Number		Date Collected Matrix	Date Prepared	Date Analyzed	QC Bate	ch ID
Method Blank			099-10	-006-18,349		06/23/06	06/24/06	060623	L02
Comment(s): -Results were e	valuated to the	MDI co	ncentratio	ne >= to the	MDL but < RL, if found, are qualified	ad with a "l" f	lan		
Parameter	Result	RL RL	MDL	DF Qual	•	Result	_	MDL	DF Qual
Acetone	ND	10	6.1	1	1,3-Dichloropropane	ND	1.0	0.30	1
Benzene	ND	0.50	0.1	i	2,2-Dichloropropane	ND ND	1.0	0.40	1
Bromobenzene	ND	1.0	0.47	1	1,1-Dichloropropene	ND	1.0	0.40	1
Bromochloromethane	ND	1.0	0.47	1	c-1,3-Dichloropropene	ND	0.50	0.45	1
Bromodichloromethane	ND	1.0	0.27	1	t-1,3-Dichloropropene	ND	0.50	0.31	i 1
Bromoform	ND	1.0	0.62	1	Ethylbenzene	ND	1.0	0.17	1
Bromomethane	ND	10	2.9	1	2-Hexanone	ND	10	1.9	1
2-Butanone	ND	10	4.2	1	Isopropylbenzene	ND	1.0	0.24	1
n-Butylbenzene	ND	1.0	0.29	1	p-Isopropyttoluene	ND	1.0	0.21	1
sec-Butylbenzene	ND	1.0	0.21	1	Methylene Chloride	2.6	10.0	2.6	1 J
tert-Butylbenzene	ND	1.0	0.17	1	4-Methyl-2-Pentanone	ND	10	2.4	1
Carbon Disulfide	ND	10	1.0	1	Naphthalene	ND	10	0.95	1.
Carbon Tetrachloride	ND	0.50	0.42	1	n-Propyibenzene	ND	1.0	0.30	1
Chlorobenzene	ND	1.0	0.36	1	Styrene	ND	1.0	0.29	1
Chloroethane	ND	1.0	0.52	1	1,1,1,2-Tetrachloroethane	ND	1.0	0.37	1
Chloroform	ND	1.0	0.22	1	1,1,2,2-Tetrachloroethane	ND	1.0	0.37	1
Chloromethane	ND	10	1.8	1	Tetrachloroethene	ND	1.0	0.29	1
2-Chlorotoluene	ND	1.0	0.24	1	Toluene	ND	1.0	0.35	1
4-Chlorotoluene	ND	1.0	0.30	1	1,2,3-Trichlorobenzene	ND	1.0	0.39	1
Dibromochloromethane	ND	1.0	0.45	1	1,2,4-Trichlorobenzene	ND	1.0	0.35	1
1.2-Dibromo-3-Chloropropane	ND	5.0	2.5	1	1.1.1-Trichloroethane	ND	1.0	0.32	1
1.2-Dibromoethane	ND	1.0	0.81	1	1,1,2-Trichloro-1,2,2-Trifluoroet	hane ND	10	0.54	1
Dibromomethane	ND	1.0	0.42	1	1,1,2-Trichloroethane	ND	1.0	0.54	1
1,2-Dichlorobenzene	ND ·	1.0	0.24	1	Trichloroethene	ND	1.0	0.30	1
1,3-Dichlorobenzene	ND	1.0	0.38	1	Trichlorofluoromethane	ND	10	0.36	1
1,4-Dichlorobenzene	ND	1.0	0.30	1	1,2,3-Trichloropropane	ND	5.0	2.3	1
Dichlorodifluoromethane	ND	1.0	0.27	1	1,2,4-Trimethylbenzene	ND	1.0	0.26	1
1,1-Dichloroethane	ND	1.0	0.53	1	1,3,5-Trimethylbenzene	ND	1.0	0.19	1
1,2-Dichloroethane	ND	0.50	0.22	1	Vinyl Acetate	ND	10	3.2	1
1,1-Dichloroethene	ND	1.0	0.31	1	Vinyl Chloride	ND	0.50	0.33	1
c-1,2-Dichloroethene	ND	1.0	0.35	1	p/m-Xylene	ND	1.0	0.38	, 1
t-1,2-Dichloroethene	ND	1.0	0.29	1	o-Xylene	ND	1.0	0.21	1
1,2-Dichloropropane	ND	1.0	0.28	1	Methyl-t-Butyl Ether (MTBE)	ND	1.0	0.29	1
Surrogates:	REC (%)	Control	<u>Limits</u>	<u>Qual</u>	Surrogates:	<u>REC (%</u>	(a) Control	<u>Limits</u>	<u>Qual</u>
Dibromofluoromethane	112	74-140			1,2-Dichloroethane-d4	108	74-146		
Toluene-d8	100	88-112			1,4-Bromofluorobenzene	81	74-110		

RL - Reporting Limit ,

DF - Dilution Factor



Calscience Environmental Laboratories, Inc.

EPA 8260B Tentatively Identified Compound List

Work Order	CEL Sample	Client ID	Q	Compound	CAS NUMBER	RT	On Column Conc. I	Estimated Conc. ug/L
06-06-1183	1	LTB061906		No TICs Found				
06-06-1183	2	MW-6		No TICs Found				
06-06-1183	3	MW-5		No TICs Found				
06-06-1183	4	MW-4		No TICs Found				
06-06-1183	5	MW-4D		Pentane	109-66-0	2.69	5.9	5.9
06-06-1183	6	LFB061906		No TICs Found				

Q Qualifier RT Retention Time





Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Date Received:

Work Order No: Preparation:

Method:

06/19/06

06-06-1183 EPA 5030B

SRL 524M-TCP

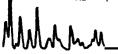
Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Page 1 of 1

Client Sample Number			Lab Sample Number		Matrix	Date Prepared	Date Analyzed	QC Batch ID
MW-6			1183-2	06/19/06	Aqueous	06/21/06	06/22/06	060621L02
<u>Parameter</u>	Result	<u>RL</u>	<u>MDL</u>	<u>DF</u>	Qual	<u>Units</u>		
1,2,3-Trichloropropane	0.19	0.02	0.0017	5	В	ug/L		
MW-5		06-06-	1183-3	06/19/06	Aqueous	06/21/06	06/22/08	060621L02
<u>Parameter</u>	Result	<u>RL</u>	MDL.	<u>DF</u>	Qual	<u>Units</u>		
1,2,3-Trichloropropane	0.14	0.01	0.0033	2		ug/L		
MW4*85		-06-06-	1183-4	1- 06/19/06	Aqueous	06/21/06	06/22/06	0606211.02
<u>Parameter</u>	Result	<u>RL</u>	MDL	DF	Qual	<u>Units</u>		
1,2,3-Trichloropropane	0.028	0.005	0.0017	1		ug/L		
MW-4D		06-06-	1183-5 👫 🕸	06/19/06	Aqueous	06/21/06	06/22/06	060621L02
<u>Parameter</u>	Result	<u>RL</u>	MDL	<u>DF</u>	<u>Qual</u>	<u>Units</u>		
1,2,3-Trichloropropane	0.027	0.005	0.0017	1	•	ug/L		
Method Blank		099-10	-022-240	NA	Aqueous	- 06/21/06 ·	06/22/06	060621L02
Comment(s): -Results were en	valuated to the l Result	MDL, concentration	ns >= to the M MDL	DL but < RL, if			flag.	
	<u> </u>	0.0050		_	<u>Qual</u>	<u>Units</u>		
1,2,3-Trichloropropane	ND		0.0017	1	***	ug/L		n ang ang ang ang ang ang ang ang ang an
Method Blank		MARIN CHINARASHARI CARANTATA IN SIN	-022-242	And the Control of th	Aqueous	06/26/06	06/26/06	060626L01
Comment(s): -Results were en	valuated to the I Result	MDL, concentration <u>RL</u>	ns >= to the Mi <u>MDL</u>	DL but < RL, if t <u>DF</u>	found, are qua Qual	lified with a "J" <u>Units</u>	flag.	
1,2,3-Trichloropropane	0.0026	0.0050	0.0017	1	J	ug/L		

RL - Reporting Limit

DF - Dilution Factor



alscience nvironmental aboratories, Inc.

Analytical Report



Tetra Tech, Inc.

Date Received:

06/19/06

3475 East Foothill Blvd., Suite 300

Work Order No:

06-06-1183

Pasadena, CA 91107-6024

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

0.050

0.01

ND

7.36

0.042

Page 1 of 2

		L	ab Sample	Number	Date				
Client Sample Number					Collected	Matri	X	_	
MW-6	4.00	$\{y,y\}$	06-06-118	3-2	06/19/06	Aqueo	us		*, A , 3 2 2 3
Comment(s): (1) Results were		ADI conce	entrations >	= to the	MDI but < F	2) if found	are qualified with	a " l" flag	20 ACC AND A STATE OF THE STATE
Parameter	Result	RL	MDL	DF	Qual	Units	Date Prepared	Date Analyzed	Method
Chromium, Hexavalent	2.9	0.2	0.0050	1	В	ug/L	N/A	06/19/06	EPA 218.6
Chloride	38	5	0.055	5		mg/L	N/A	06/20/06	EPA 300.0
Nitrite (as N) (1)	ND	0.10	0.015	1		mg/L	N/A	06/20/06	EPA 300.0
Nitrate (as N)	13	0.50	0.028	5		mg/L	N/A	06/20/06	EPA 300.0
Sulfate	74	10	0.069	10		mg/L	N/A	06/20/06	EPA 300.0
Perchlorate (1)	0.56	2.0	0.43	1	j	ug/L	N/A	06/23/06	EPA 314.0
Sulfide, Total (1)	ND	0.050	0.042	1		mg/L	06/22/06	06/22/06	EPA 376.2
Dissolved Oxygen	8.17	0.01		1		mg/L	N/A	06/19/06	SM 4500-O G

MW-5			06-06-118	3-3	06/19/06	S Aque	ous ,		in the second
Comment(s): (1) Results wer	e evaluated to the N	MDL, conce	entrations >	= to the	MDL but	RL, if foun	d, are qualified with	h a "J" flag.	
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	MDL	<u>DF</u>	Qual	<u>Units</u>	Date Prepared	Date Analyzed	<u>Method</u>
Chromium, Hexavalent	1.7	0.2	0.0050	1	В	ug/L	N/A	06/19/06	EPA 218.6
Chloride	37	5	0.055	5		mg/L	N/A	06/20/06	EPA 300.0
Nitrite (as N) (1)	ND	0.10	0.015	1		mg/L	N/A	06/20/06	EPA 300.0
Nitrate (as N)	13	0.50	0.028	5		mg/L	N/A	06/20/06	EPA 300.0
Sulfate	75	10	0.069	10		mg/L	N/A	06/20/06	EPA 300.0
Perchlorate (1)	0.99	2.0	0.43	1	J	ug/L	N/A	06/23/06	EPA 314.0

MW4		i Herita	06-06-118	3-4,	06/19/0	6 Aque	ous .		
Comment(s): (1) Results were	evaluated to the N	1DL, conce	entrations >	= to the	MDL but	< RL, if foun	d, are qualified with	n a "J" flag.	
<u>Parameter</u>	Result	<u>RL</u>	MDL	<u>DF</u>	<u>Qual</u>	<u>Units</u>	Date Prepared	Date Analyzed	<u>Method</u>
Chromium, Hexavalent	1.5	0.2	0.0050	1	В	ug/L	N/A	06/19/06	EPA 218.6

<u>Parameter</u>	<u>Result</u>	<u>KL</u>	MUL	<u>DF</u>	<u>Qual</u>	<u>Units</u>	Date Prepared	Date Analyzed	Method	
Chromium, Hexavalent	1.5	0.2	0.0050	1	В	ug/L	N/A	06/19/06	EPA 218.6	
Chloride	40	10	0.055	10		mg/L	N/A	06/20/06	EPA 300.0	
Nitrite (as N) (1)	ND	0.10	0.015	1		mg/L	N/A	06/20/06	EPA 300.0	
Nitrate (as N)	16	0.50	0.028	5		mg/L	N/A	06/20/06	EPA 300.0	
Sulfate	72	10	0.069	10		mg/L	N/A	06/20/06	EPA 300.0	
Perchlorate (1)	0.58	2.0	0.43	1	J	ug/L	N/A	06/23/06	EPA 314.0	
Sulfide, Total (1)	ND	0.050	0.042	1		mg/L	06/22/06	06/22/06	EPA 376.2	
Dissolved Oxygen	7.89	0.01		1		mg/L	N/A	06/19/06	SM 4500-O G	

RL - Reporting Limit ,

DF - Dilution Factor

Qual - Qualifiers

Sulfide, Total (1)

Dissolved Oxygen

mg/L

mg/L

N/A

06/22/06

06/22/06

06/19/06

EPA 376.2

SM 4500-O G





Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Date Received:

06/19/06

Work Order No:

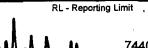
06-06-1183

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Page 2 of 2

Client Sample Number			ab Sample	Numbe	r Date Collected	_I Matri	×		· .
MW-4D			06-06-118	3-5	06/19/06	Aqueo	nus (**) 1	Page 1	
Comment(s): (1) Results wer	re evaluated to the N	1DL, conce	entrations >	= to the	MDL but <	RL, if found	I, are qualified with	h a "J" flag.	
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>MDL</u>	<u>DF</u>	Qual	<u>Units</u>	Date Prepared	Date Analyzed	<u>Method</u>
Chromium, Hexavalent	1.5	0.2	0.0050	1	В	ug/L	N/A	06/19/06	EPA 218.6
Chloride	44	5	0.055	5		mg/L	N/A	06/20/06	EPA 300.0
Nitrite (as N) (1)	ND	0.10	0.015	1		mg/L	N/A	06/20/06	EPA 300.0
Nitrate (as N)	12	0.50	0.028	5		mg/L	N/A	06/20/06	EPA 300.0
Sulfate	73	10	0.069	10		mg/L	N/A	06/20/06	EPA 300.0
Perchlorate (1)	0.71	2.0	0.43	1	· J	ug/L	N/A	06/23/06	EPA 314.0
Sulfide, Total (1)	ND	0.050	0.042	1		mg/L	06/22/06	06/22/06	EPA 376.2
Dissolved Oxygen	7.82	0.01		1		mg/L	N/A	06/19/06	SM 4500-O G

MethodiBlank					ŅĀ	Aque	ous 🗀 🕌	. 14		
Comment(s): (1) Results were	evaluated to the M	MDL, conce	entrations >	= to the	MDL but	RL, if found	d, are qualified with	n a "J" flag.		
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	MDL	<u>DE</u>	Qual	<u>Units</u>	Date Prepared	Date Analyzed	<u>Method</u>	
Chromium, Hexavalent (1)	0.025	0.20	0.0050	1	J	ug/L	N/A	06/19/06	EPA 218.6	
Chloride (1)	ND	1.0	0.055	1		mg/L	N/A	06/20/06	EPA 300.0	
Nitrite (as N) (1)	ND	0.10	0.015	1		mg/L	N/A	06/20/06	EPA 300.0	
Nitrate (as N) (1)	ND	0.10	0.028	1		mg/L	N/A	06/20/06	EPA 300.0	
Sulfate (1)	ND	1.0	0.069	1		mg/L	N/A	06/20/06	EPA 300.0	
Perchlorate (1)	ND	2.0	0.43	1		ug/L	N/A	06/23/06	EPA 314.0	
Sulfide, Total (1)	ND	0.050	0.042	1		mg/L	06/22/06	06/22/06	EPA 376.2	



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Quality Control - Spike/Spike Duplicate



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: 06/19/06 06-06-1183 EPA 3005A Filt. EPA 6010B

Project BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID	Matrix	Instrument	Date Prepared	Α	Date nalyzed	MS/MSD Batch Number
MW-4	Aqueous	i - ICP 3300	06/20/06	0	6/21/06	/060620S03
Parameter	MS %REC	MSD %REC	%REC CL	<u>RPD</u>	RPD CL	Qualifiers
Antimony	105	108	72-132	3 .	0-10	
Arsenic	104	108	80-140	4	0-11	
Barium	103	105	87-123	1	0-6	
Beryllium 1	100	105	89-119	4	8-0	
Cadmium	102	104	82-124	2	0-7	
Chromium	103	105	86-122	2	0-8	
Cobatt	99	105	83-125	6	0-7	
Copper	70	72	78-126	3	0-7	3
Lead	100	103	84-120	3	0-7	
Molybdenum	103	106	78-126	3	0-7	
Nickel	96	102	84-120	6	0-7	
Selenium	104	106	79-127	2	0-9	
Silver	104	105	86-128	1	0-7	
Thallium	. 94	98	79-121	4	0-8	
Vanadium	103	105	88-118	2	0-7	
Zinc	98	105	89-131	6	0-8	

MAM_



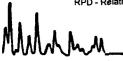


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Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: 06/19/06 06-06-1183 EPA 3005A Filt. EPA 6010B

Project BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID	Matrix	Instrument	Date Prepared	,	Date Analyzed	MS/MSD Batch Number
WW4	Aqueou	s ICP 3300	06/20/06	1	06/21/06	060620503
<u>Parameter</u>	MS %REC	MSD %REC	%REC CL	RPD	RPD CL	Qualifiers
Calcium	4X	4X	77-113	4X	0-11	Q
Magnesium	4X	4X	56-140	4X	0-11	Q
Potassium	102	102	83-131	0	0-7	
Sodium	4X	. 4X	73-127	4X	0-9	Q



RPD - Relative Percent Difference , CL - Control Limit

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Quality Control - Spike/Spike Duplicate

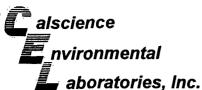


Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: 06/19/06 06-06-1183 EPA 3005A Filt. EPA 200.8

Project BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID	Matrix	Instrument	Date Prepared		Date alyzed	MS/MSD Batch Number
MW4	Aqueou	s ICP/MS A	06/21/06	Ó	V21/06	060621801
<u>Parameter</u>	MS %REC	MSD %REC	%REC CL	RPD	RPD CL	Qualifiers
Iron Manganese	142 100	134 100	80-120 80-120	6	0-20 0-20	3

hhm_





etra Tech Inc

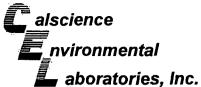
Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: 06/19/06 06-06-1183 EPA 7470A Filt. EPA 7470A

Project BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID	Matrix	Instrument	Date Prepared	Date Analyzed	MS/MSD Batch Number
MW4 (1) (1)	Aqueou	s 🤲 Mercury	06/20/06	06/20/06	060620802
<u>Parameter</u>	MS %REC	MSD %REC	%REC CL	RPD RP	D CL Qualifiers
Mercury	83	83	71-134	0 0-	-14

Mulham

RPD - Relative Percent Difference , CL - Control Lin





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Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: 06/19/06 06-06-1183 EPA 3520B EPA 8270C(M) Isotope Dilution

Project BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID	Matrix	Instrument	Date Prepared	i An	alyzed	MS/MSD Batch Number
	Aqueou	GC/MS-J	06/22/06		/28/06	060622S01D
<u>Parameter</u>	MS %REC	MSD %REC	%REC CL	RPD	RPD CL	Qualifiers
1,4-Dioxane	83	85	50-130	3	0-20	

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Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received: Work Order No: Preparation: Method: 06/19/06 06-06-1183 EPA 3520B EPA 1625CM

Project BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID	Matrix	Instrument	Date Prepared		Date nalyzed	MS/MSD Batch Number
MW4	Aqueou	s GCMSH	06/23/06	3 (\$) (6/26/06	060623S09A
<u>Parameter</u>	MS %REC	MSD %REC	%REC CL	RPD	RPD CL	Qualifiers
N-Nitrosodimethylamine	118	99	50-130	17	0-20	

All RPD - Relati

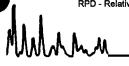
RPD - Relative Percent Difference , CL - Control I

Quality Control - Spike/Spike Duplicate



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: 06/19/06 06-06-1183 EPA 5030B EPA 8260B

Quality Control Sample ID	Matrix	Instrument	Date Prepared		Date Analyzed	MS/MSD Batch Number	
MW4	Aqueous	s⊭ ∴GC/MSW	06/22/06	06/22/06		060622802	
<u>Parameter</u>	MS %REC	MSD %REC	%REC CL	RPD	RPD CL	Qualifiers	
Benzene	98	97	88-118	1	0-7		
Carbon Tetrachloride	101	108	67-145	7	0-11		
Chlorobenzene	104	100	88-118	4	0-7		
1,2-Dichlorobenzene	103	101	86-116	2	0-8		
1,1-Dichloroethene	93	96	70-130	4	0-25		
Toluene	102	98	87-123	5	0-8		
Trichloroethene	86	86	79-127	0	0-10		
Vinyl Chloride	93	106	69-129	14	0-13	4	
Methyl-t-Butyl Ether (MTBE)	88	101	71-131	14	0-13	4	
Tert-Butyl Alcohol (TBA)	79	95	36-168	18	0-45		
Diisopropyl Ether (DIPE)	93	97	81-123	4	0-9	•	
Ethyl-t-Butyl Ether (ETBE)	86	94	72-126	8	0-12		
Tert-Amyl-Methyl Ether (TAME)	99	101	72-126	2	0-12		
Ethanol	115	87	53-149	28	0-31		



Quality Control - Spike/Spike Duplicate



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: 06/19/06 06-06-1183 EPA 5030B EPA 8260B

Project BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID	Matrix	Instrument	Date ument Prepared		Date Analyzed	MS/MSD Batch Number
06-08-1352-1	Aqueous	GC/MS W	06/23/06		06/24/06	060623S02
<u>Parameter</u>	MS %REC	MSD %REC	%REC CL	RPD	RPD CL	Qualifiers
•						
Benzene	92	89	88-118	3	0-7	
Carbon Tetrachloride	93	94	67-145	1	0-11	
Chlorobenzene	98	97	88-118	1	0-7	
1,2-Dichlorobenzene	98	97	86-116	1	8-0	
1,1-Dichloroethene	83	82	70-130	1	0-25	
Toluene	92	89	87-123	3	0-8	
Trichloroethene	94	90	79-127	4	0-10	
Vinyl Chloride	88	87	69-129	1	0-13	
Methyl-t-Butyl Ether (MTBE)	0	0	71-131	8	0-13	3
Tert-Butyl Alcohol (TBA)	84	84	36-168	0	0-45	
Diisopropyl Ether (DIPE)	89	89	81-123	1	0-9	
Ethyl-t-Butyl Ether (ETBE)	76	77	72-126	1	0-12	
Tert-Amyl-Methyl Ether (TAME)	90	89	72-126	1	0-12	
Ethanol	137	134	53-149	2	0-31	

RPD - Rela

Quality Control - Spike/Spike Duplicate



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: 06/19/06 06-06-1183 EPA 5030B SRL 524M-TCP

Project BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID	Matrix	Matrix Instrument		d A	Date nalyzed	MS/MSD Batch Number	
MW-4	Aqueou	s GC/MS M	06/21/06	# # C	6/22/06	060621302	
<u>Parameter</u>	MS %REC	MSD %REC	%REC CL	<u>RPD</u>	RPD CL	Qualifiers	
1,2,3-Trichloropropane 1,4-Dioxane	66 90	66 89	80-120 80-120	0 2	0-20 · 0-20	3	

MMM



Quality Control - Spike/Spike Duplicate



aboratories, Inc.

Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation:

Method:

06/19/06 06-06-1183 EPA 5030B SRL 524M-TCP

Project BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID	Matrix	Instrument	1 (cpaica		nalyzed	MS/MSD Batch Number	
06-06-1520-1		s ∵GC/MS M≀	06/26/06	.	6/26/06	060626S01	
<u>Parameter</u>	MS %REC	MSD %REC	%REC CL	<u>RPD</u>	RPD CL	Qualifiers	
1,2,3-Trichloropropane 1,4-Dioxane	94 100	124 100	80-120 80-120	28 0	0-20 0-20	4,3	

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RPD - Relative Percent Difference , CL - Control Limit

Quality Control - Spike/Spike Duplicate



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Date Received: Work Order No:

N/A 06-06-1183

Matrix: Aqueous#			74				100	7.		
<u>Parameter</u>	<u>Method</u>	Quality Control Sample ID	<u>Date</u> <u>Analyzed</u>	<u>Date</u> <u>Extracted</u>	MS% REC	MSD % REC	%REC CL	<u>RPD</u>	RPD CL	Qualifiers
Chloride	EPA 300.0	MW-4	06/20/06	N/A	95	96	56-134	1	0-3	
Nitrite (as N)	EPA 300.0	MW-4	06/20/06	N/A	99	95	68-122	4	0-8	
Nitrate (as N)	EPA 300.0	MW-4	06/20/06	N/A	93	96	58-142	4	0-6	
Sulfate	EPA 300.0	MW-4	06/20/06	N/A	104	102	49-133	2	0-3	
Chromium, Hexavalent	EPA 218.6	MW-4	06/19/06	N/A	103	102	85-121	0	0-4	
Perchlorate	EPA 314.0	MW-4	06/23/06	N/A	97	97	80-120	0	0-15	



Quality Control - Duplicate

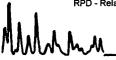


Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No:

1N/*F*

06-06-1183

Matrix: Aqueous .			i Kar				7.	
<u>Parameter</u>	Method	QC Sample ID	Date Analyzed	Sample Conc	DUP Conc	RPD	RPD CL	Qualifiers
Dissolved Oxygen Sulfide, Total	SM 4500-O G EPA 376.2	MW-4 MW-4	06/19/06 06/22/06	7.89 ND	7.70 ND	2 NA	0-25 0-25	



alscience nvironmental Quality Control - Laboratory Control Sample aboratories, Inc.



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method:

06-06-1183 EPA 3005A Filt. EPA 6010B

Quality Control Sample ID	Matrix	Instrum	ent Date Analyzed	Lab File	ID LCS B	atch Number
097-01-003-6,215	Aqueous	. ICP 33	06/21/06	0606204	.03 6	0620L03F
<u>Parameter</u>	<u>9</u>	Conc Added	Conc Recovered	LCS %Rec	%Rec CL	<u>Qualifiers</u>
Antimony		0.500	0.528	106	80-120	
Arsenic		0.500	0.522	104	80-120	
Barium		0.500	0.533	107	80-120	
Beryllium		0.500	0.512	102	80-120	
Cadmium	•	0.500	0.535	107	80-120	
Chromium		0.500	0.525	105	80-120	
Cobalt		0.500	0.544	109	80-120	
Copper		0.500	0.447	89	80-120	
Lead		0.500	0.532	106	80-120	
Molybdenum		0.500	0.530	106	80-120	
Nickel		0.500	0.543	109	80-120	
Selenium		0.500	0.493	99	80-120	
Silver		0.250	0.263	105	80-120	
Thallium		0.500	0.489	98	80-120	
Vanadium		0.500	0.512	102	80-120	
Zinc		0.500	0.529	106	80-120	



alscience nvironmental Quality Control - Laboratory Control Sample aboratories, Inc.



Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

Work Order No:

Preparation: Method:

N/A

06-06-1183

EPA 3005A Filt.

EPA 6010B

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID	Matrix	Instrume	ent Date Analyze	d Lab File	ID LCS E	Batch Number
097-01-003-6;215	Aqueous	- CP 330	06/21/06	-0606204	-03 🕌 06	0620L03F
<u>Parameter</u>		Conc Added	Conc Recovered	LCS %Rec	%Rec CL	Qualifiers
Calcium		0.500	0.516	103	80-120	
Magnesium		0.500	0.511	102	80-120	
Potassium		5.00	5.12	102	80-120	
Sodium		5.00	5.20	104	80-120	

RPD - Relative Percent Difference,

CL - Control Limit

Quality Control - LCS/LCS Duplicate



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Date Received: Work Order No: Preparation: Method: N/A 06-06-1183 EPA 3005A Filt. EPA 200.8

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID		Instrument	Date Prepared	Date Analyzed	LCS/LCSD Bate Number	ch
099_10_008-736	Aqueous:	ICP/MS A	06/21/08	06/21/06	-060621L01	
Parameter	LCS %RE	C LCSD %	REC %REC	CCL RPD	RPD CL	Qualifiers
Iron	102	102	85-1	15 0	0-20	
Manganese	98	98	85-1	15 0	0-20	

RPD - Relative Percent Difference ,

alscience nvironmental Quality Control - Laboratory Control Sample aboratories, Inc.



Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Date Received:

Work Order No:

Preparation: Method:

06-06-1183

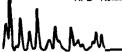
N/A

EPA 7470A Filt.

EPA 7470A

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID	Matrix-	Instrument	Date Analyzed	Lab File I	LCS	Batch Number
2099-04-008-2/54-1	Aqueous 1	Mercury	06/21/06	060620-1-02-1	icp 👫 🖑	060620L02F
Parameter	Co	nc Added	Conc Recovered	LCS %Rec	%Rec CL	Qualifiers
Mercury	(0.0100	0.00970	97	90-122	



PD - Relative Percent Difference,

CL - Control Limit

Quality Control - LCS/LCS Duplicate



Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Date Received:

N/A

Work Order No:

06-06-1183

Preparation:

EPA 3520B

Method:

EPA 8270C(M) Isotope Dilution

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID	Matrix I	nstrument	Date Prepared	Date Analyzed	LCS/LCSD Bate Number	ch
099-09-004-597.	queous .	GC/MS J	06/22/06	06/27/06	060622L01D	
<u>Parameter</u>	LCS %REC	LCSD %	REC %REC	CL RPD	RPD CL	Qualifiers
1,4-Dioxane	83	88	50-13	30 5	0-20	

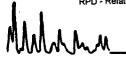
Mhha_





Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: N/A 06-06-1183 EPA 3520B EPA 1625CM

Quality Control Sample ID	Matrix Ir	nstrument	Date Prepared	Date Analyzed	LCS/LCSD Ba Number	tch
099-07-027-251	Aqueous 🔭 🐮 🤆 🤇	SC/MS H	06/23/06 #	06/27/06	060623L09	
Parameter	LCS %REC	LCSD %	REC %REC	C CL RP	D RPD CL	Qualifiers
N-Nitrosodimethylamine	78	80	50-	130 4	0-20	







Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: N/A 06-06-1183 EPA 5030B EPA 8260B

Quality Control Sample ID	encontract and a second	Instrument		Date Analyzed 06/22/06	LCS/LCSD Bate Number	
USS IV WO IO SEA	Aqueous	goma ve	U0/22/U0	0/2/200	VOUGEZEUZ	1. T. A.
<u>Parameter</u>	LCS %REC	LCSD %R	EC %REC CI	RPD	RPD CL	Qualifiers
Benzene	97	97	84-120	0	0-8	
Carbon Tetrachloride	104	110	63-147	6	0-10	
Chlorobenzene	100	103	89-119	3	0-7	
1,2-Dichlorobenzene	102	103	89-119	1	0-9	
1,1-Dichloroethene	91	96	77-125	6	0-16	
Toluene	99	98	83-125	0	0-9	
Trichloroethene	100	101	89-119	1	0-8	
Vinyl Chloride	104	107	63-135	3	0-13	
Methyl-t-Butyl Ether (MTBE)	93	102	82-118	9	0-13	
Tert-Butyl Alcohol (TBA)	90	97	46-154	7	0-32	
Diisopropyl Ether (DIPE)	92	99	81-123	7	0-11	
Ethyl-t-Butyl Ether (ETBE)	89	98	74-122	10	0-12	
Tert-Amyl-Methyl Ether (TAME)	99	104	76-124	6	0-10	
Ethanol	95	96	60-138	2	0-32	



Quality Control - LCS/LCS Duplicate



Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

Work Order No:

Preparation:

Method:

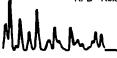
N/A

06-06-1183

EPA 5030B

EPA 8260B

Quality Control Sample ID	Matrix Instrument		Date Prepared 06/23/06	Date Analyzed	LCS/LCSD Bate Number 060623L02	th
		Zame.	ourojoo	002900		AND THE STREET
Parameter	LCS %REC	LCSD %F	EC %REC	CL RPD	RPD CL	Qualifiers
Benz <i>e</i> ne	96	94	84 -1	20 3	0-8	
Carbon Tetrachloride	106	103	63-1	147 3	0-10	
Chlorobenzene	101	99	89-1	119 1	0-7	•
1,2-Dichlorobenzene	100	99	89-1	119 1	0-9	
1,1-Dichloroethene	94	92	77-1	25 3	0-16	
Toluene	98	97	83-1	25 0	0-9	
Trichloroethene	99	96	89-1	19 3	0-8	
Vinyl Chloride	105	92	63-1	35 13	0-13	
Methyl-t-Butyl Ether (MTBE)	92	82	82-1	118 11	0-13	
Tert-Butyl Alcohol (TBA)	82	. 78	46 -1	54 5	0-32	
Diisopropyl Ether (DIPE)	94	91	81-1	23 3	0-11	
Ethyl-t-Butyl Ether (ETBE)	89	85	74-1	22 4	0-12	
Tert-Amyl-Methyl Ether (TAME)	97	96	76- 1	24 2	0-10	
Ethanol A	81	112	60 -1	38 32	0-32	







Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Date Received:

Work Order No:

Preparation:

Method:

N/A

06-06-1183

EPA 5030B

SRL 524M-TCP

Quality Control Sample ID	Matrix Ir	nstrument	Date Prepared	Date Analyzed	LCS/LCSD Batch Number	· ·
099-10-022-240	Aqueous G	C/MS M	06/21/06	06/21/06	* 060621L02	
<u>Parameter</u>	LCS %REC	LCSD %R	EC %REC	CL RPD	RPD CL	Qualifiers
1,2,3-Trichloropropane	86	91	80-12	20 6	0-20	
1,4-Dioxane	83	87	80-12	20 4	0-20	







Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

Work Order No:

Preparation:

Method:

N/A

06-06-1183

EPA 5030B SRL 524M-TCP

Quality Control Sample ID		strument	Date Prepared	Date Analyzed	LCS/LCSD Batc Number 060626L01	
099-10-022-242 Parameter	LCS %REC	C/MS M LCSD %R	AC CHAIRMAN AND AND AND AND AND AND AND AND AND A	06/26/06	060626E01	Qualifiers
1,2,3-Trichloropropane 1,4-Dioxane	107 94	115 94	80-120 80-120	7	0-20 0-20	



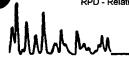


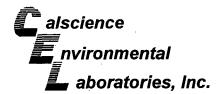
Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received:

N/A

Work Order No: 06-06-1183

Matrix Aqueous				r ilk						
<u>Parameter</u>	<u>Method</u>	Quality Control Sample ID	<u>Date</u> Extracted	<u>Date</u> <u>Analyzed</u>	LCS % REC	LCSD % REC	%REC CL	<u>RPD</u>	RPD CL	Qual
Chloride	EPA 300.0	099-05-118-3,422	N/A	06/20/06	94	95	81-111	1	0-5	
Nitrite (as N)	EPA 300.0	099-05-118-3,422	N/A	06/20/06	92	94	73-115	1	0-26	
Nitrate (as N)	EPA 300.0	099-05-118-3,422	N/A	06/20/06	97	97	87-111	0	0-12	
Sulfate	EPA 300.0	099-05-118-3,422	N/A	06/20/06	99	97	89-107	2	0-13	
Chromium, Hexavalent	EPA 218.6	099-05-124-485	N/A	06/19/06	99	99	95-107	0	0-20	
Perchlorate	EPA 314.0	099-05-203-423	N/A	06/23/06	107	104	85-115	3	0-15	





Glossary of Terms and Qualifiers



Work Order Number: 06-06-1183

Qualifier	<u>Definition</u>
*	See applicable analysis comment.
1	Surrogate compound recovery was out of control due to a required sample dilution, therefore, the sample data was reported without further clarification.
2	Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification.
3	Recovery of the Matrix Spike or Matrix Spike Duplicate compound was out of control due to matrix interference. The associated LCS and/or LCSD was in control and, therefore, the sample data was reported without further clarification.
4	The MS/MSD RPD was out of control due to matrix interference. The LCS/LCSD RPD was in control and, therefore, the sample data was reported without further clarification.
5	The PDS/PDSD associated with this batch of samples was out of control due to a matrix interference effect. The associated batch LCS/LCSD was in control and, hence, the associated sample data was reported with no further corrective action required.
Α	Result is the average of all dilutions, as defined by the method.
В	Analyte was present in the associated method blank.
С	Analyte presence was not confirmed on primary column.
Е	Concentration exceeds the calibration range.
Н	Sample received and/or analyzed past the recommended holding time.
J	Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated.
Ν .	Nontarget Analyte.
ND	Parameter not detected at the indicated reporting limit.
Q	Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater.
υ	Undetected at the laboratory method detection limit.
Χ	% Recovery and/or RPD out-of-range.
Z	Analyte presence was not confirmed by second column or GC/MS analysis.

	TETRA TECH, INC.
	3475 E. FOOTHILL BLVD. PASADENA, CALIFORNIA 91107 TELEPHONE (626) 351-4664 FAX (626)351-5291
	PASADENA, CALIFORNIA 91107
	TELEPHONE (626) 351-4664
· · · · · · · ·	FAX (626)351-5291

SHIPPED TO: CALSCIENCE
7440 Uneder War

CHAIN OF CUSTODY RECORD

DATE 6/19/66 PAGE 1 OF 1

SAMPLE NO. DATE TIME STELL SELVE PRESERVATIVES HOLL MATRIX TYPE 8-SOL. WAS GO 2 MAY - CONTAINER G-GLAS BOTTLEWA TYPE 88-STRINLESS STELL SLEEVE MATRIX TYPE 8-SOL. WAS GO 2 MAY - CONTAINER G-GLAS BOTTLEWA TYPE 88-STRINLESS STELL SLEEVE MATRIX TY	FAX (626)351-5291	1					<u>ا با7</u>	Σ	<u>/ L</u>	<i>I</i> D (9	100	71				DAT	E _	2(1	7	90	PA	GE —	_OF _	
STANDARD STANDA	CLIENT LOCKHEED MA	ARTIN CO	>RP			2	EXT	rac'	TION/	ANAL	Sic	AL ME	ТНО	DS			•								
TABLE NO. DATE TIME SAMPLERS (SIGNATURE) SAMPLE NO. DATE TIME SAMPLE NO. DATE SAMPLE NO. DATE TIME SAMPLE NO. DATE TIME SAMPLE NO. DATE SAMPLE	PROJECT NAME: 2006 E	30U			**	म	2			,	35	,	`										TURN-	ROUND	TIME
SAMPLE NO. DATE TIME SAMPLE NO. DATE TIME X X X X X X X X X X X X X X X X X X X	Graundwater Monitoria	ng (PACH	IELLS)	SS	家	14	2					٥	5	80				·	}	١	22	1			
SAMPLE NO. DATE TIME SAMPLE NO. DATE TIME X X X X X X X X X X X X X X X X X X X	TASK MANAGER: Neil S	hu.K.la.		300	\$ 21 20 20 20 20 20 20 20 20 20 20 20 20 20	ags Sec	104	<u>2</u>	450		11.75 11.75	× 00	22	18 N	\$						AINE	. [NS
SAMPLE NO. DATE TIME SUBJECT TIME SECRETATIVES HOLL NOR (NONE REQUIRED) MATERIAL TYPE S. DOL. W. WATER S. SUDDE SIGNATURE COMPANY LOTAL NUMBER OF CONTAINER SUBJECT SIGNATURE COMPANY LOTAL NUMBER SIGNATURE COMPANY LOTAL NUMBER SIGNATURE SIGNATURE COMPANY LOTAL NUMBER SIGNATURE SIGNATURE SIGNATURE SIGNATURE COMPANY LOTAL NUMBER SIGNATURE SIGNATU	TC#: 17597-3.1A	P17653-E	602	ώy M	EP	がな	÷%	20	√ 16	24 50	86	3.5	$\mathcal{O}_{\mathbf{Z}}$	E	20				Ψ,		8				1
LTB061966 6/19/66 0805 X X X X X X X X X X X X X X X X X X X	SAMPLERS (SIGNAPHIRE)	R,		Ss EP	23	6.22	A Zi	- Plok	W ER	242 30	Solved 200	STAN E	solves	नार्टा <i>न</i>	trate/			RIX TYPE	SERVATIN	TAINER	BER OF				
LTB061966 6/19/66 0805 X X X X X X X X X X X X X X X X X X X	SAMPLE NO.	DATE	TIME	ğξ	(52	TIT EPA	四至	1,4 6,9	N	55 m	PIS EPA	30	3	S	1 2 1 1	.	.	MAT	PRE	8	N S				
MW-6 D945 X X X X X X X X X X X X X X X X X X X				X	X	X	X	X	X		X	X	X	X	X			·							
MW-5 1205 X X X X X X X X X X X X X X X X X X X	LT B061966	6/19/06	0805	X											·			W	Yes	بل	2				
MW-5 1205 X X X X X X X X X X X X X X X X X X X	MW-6		0945	X	X	X	X	X	X	X	X	X	K	X	X			W	Yes	9/0	13				
MATRIX TYPE: 8 - SOIL W. WATER S. SLUDGE RELINQUISHED BY RELINQUISHED BY LONG AND	MW-5		1205	X	X	Χ	X	X	人	X	X	×	又	X	X			W			13				
MATRIX TYPE: S. SOIL W. YES G. 2 MATRIX	MW-9		1329	K	X	X	Х	又	X	又	X	X	X	X	K				, , , , , , , , , , , , , , , ,	11	.02				•
MATRIX TYPE: 8 - SOIL. WYS G 2 MATRIX TYPE: 8 - SOIL. NR (NONE REQUIRED) TEMPERATURE BLANK WES NO TETRA TECH, INC. DATE TIME METHOD OF SHIPMENT METHOD O			1339	X	1	Z.	X	X	义	人	X	X	X	X	X			W	403	6/8	39		MS	us1)
MATRIX TYPE: \$ - SOIL CONTAINER G - GLASS BOTTLEVOA TYPE: SB - STAINLESS STEEL SLEEVE NR (NONE REQUIRED) RELINQUISHED BY RECEIVED BY LOUISI DO MACAS RECEIVED BY RELINQUISHED BY SIGNATURE COMPANY COMPANY COMPANY DATE TIME A FIME A FIM		\bigvee	1405	X					·									W	Yes	ઉ	2				
MATRIX TYPE: \$ - SOIL CONTAINER G - GLASS BOTTLEVOA TYPE: SB - STAINLESS STEEL SLEEVE NR (NONE REQUIRED) RELINQUISHED BY RECEIVED BY LOUISI DO MACAS RECEIVED BY RELINQUISHED BY SIGNATURE COMPANY COMPANY COMPANY DATE TIME A FIME A FIM																		,							
RECEIVED BY SIGNATURE TETRA TECH, INC. 6/9/66 1579 OF CONTAINERS PATE TIME TOTAL NUMBER OF CONTAINERS PATE TIME METHOD OF SHIPMENT OF COMPANY DATE TIME SPECIAL SHIPMENT/HANDLING OR STORAGE GUYPCY OF MAIGS																									
RECEIVED BY SIGNATURE TETRA TECH, INC. 6/9/66 1579 OF CONTAINERS PATE TIME TOTAL NUMBER OF CONTAINERS PATE TIME METHOD OF SHIPMENT OF COMPANY DATE TIME SPECIAL SHIPMENT/HANDLING OR STORAGE GUYPCY OF MAIGS																									
RECEIVED BY SIGNATURE TETRA TECH, INC. 6/9/66 1579 OF CONTAINERS RECEIVED BY SIGNATURE COMPANY DATE TIME METHOD OF SHIPMENT SIGNATURE COMPANY DATE TIME SPECIAL SHIPMENT/HANDLING OR STORAGE REQUIREMENTS																									
RECEIVED BY SIGNATURE COMPANY DATE TIME METHOD OF SHIPMENT HANDLING OR STORAGE REQUIREMENTS	MATRIX TYPE: S - SOIL CONT. W - WATER SL - SLUDGE	AINER G - GLASSI YPE: SS - STAINLE P - PLASTE	BOTTLENOA ESS STEELS C	LEEVE		PRES	RVATI			IE REQI	UIRED)	•												NO
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RELINQUISHED BY SIGNATURE COMPANY DATE TIME SPECIAL SHIPMENT/HANDLING OR STORAGE REQUIREMENTS	LEONIGIDO Macas			2	1_	<u> </u>	E	-ر					, PĂ	LE	15	ME /S		_/_	pl	· •——	[1]			2	
RECEIVED BY COMPANY COMPANY DATE TIME AIRBILL NO:	Loovey do Macias				1_	(t	1		·			19 19	106	172	8	REQUI	REME	NTS	ENT/H	AND	LING	OR ST)RAGE	
	RECEIVED BY			\A\		OMPA		2	7									L NC):		 -				



WORK ORDER #:	06	- 0	Le -	1	8	3

Cooler	of
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SAMPLE REC	EIPT FORM
CLIENT: TETRA TECH	DATE: 6-19-06
TEMPERATURE - SAMPLES RECEIVED BY:	
CALSCIENCE COURIER: Chilled, cooler with temperature blank provided. Chilled, cooler without temperature blank. Chilled and placed in cooler with wet ice. Ambient and placed in cooler with wet ice. Ambient temperature. Compared to cooler with wet ice. Compared to cooler with wet ice. Cooler with wet ice.	LABORATORY (Other than Calscience Courier): ° C Temperature blank. ° C IR thermometer. Ambient temperature.
CUSTODY SEAL INTACT: Sample(s): Cooler: No (Not Intact)): Not Applicable (N/A): Initial:/
SAMPLE CONDITION: Chain-Of-Custody document(s) received with samples	
COMMENTS:	





July 05, 2006

Neil Shukla Tetra Tech. Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Calscience Work Order No.: Subject:

06-06-1241

Client Reference: **BOU Groundwater Monitoring 2006 (PAC Wells)**

/ 17653-0602

Dear Client:

Enclosed is an analytical report for the above-referenced project. The samples included in this report were received 6/20/2006 and analyzed in accordance with the attached chain-of-custody.

Unless otherwise noted, all analytical testing was accomplished in accordance with the guidelines established in our Quality Systems Manual, applicable standard operating procedures, and other related documentation. The original report of any subcontracted analysis is provided herein, and follows the standard Calscience data package. The results in this analytical report are limited to the samples tested and any reproduction thereof must be made in its entirety.

If you have any questions regarding this report, please do not hesitate to contact the undersigned.

Sincerely,

Calscience Environmental Laboratories, Inc.

Jason Torres **Project Manager**

CA-ELAP ID: 1230

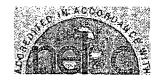
NELAP ID: 03220CA

CSDLAC ID: 10109

SCAQMD ID: 93LA0830

7440 Lincoln Way, Garden Grove, CA 92841-1427 • TEL:(714) 895-5494 • FAX: (714) 894-7501





Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

06/20/06

Work Order No:

06-06-1241

Preparation:

EPA 3005A Filt. / EPA 7470A Filt.

Method:

EPA 6010B / EPA 7470A

Units:

mg/L

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Page 1 of 2

	Lab Sample	Date		Date	Date	
Client Sample Number	Number	Collected	Matrix	Prepared	Analyzed	QC Batch ID
MVV-8	06-06-1241-2	* 06/20/06	Aqueous	06/21/06	06/22/06	060621L06F

Comment(s): -Results were evaluated to the MDL, concentrations >= to the MDL but < RL, if found, are qualified with a "J" flag.

-Mercury was analyzed on 6/22/2006 1:33:43 PM with batch 060621L04F

<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>MDL</u>	<u>DF</u>	<u>Qual</u>	<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>MDL</u>	DF Qual	
Antimony	ND	0.0150	0.00209	1		Mercury	ND	0.000500	0.0000672	1	
Arsenic	ND	0.0100	0.00308	1		Molybdenum	ND.	0.00500	0.000800	1	
Barium	0.145	0.010	0.000719	1		Nickel	0.00239	0.00500	0.00137	1 J,B	
Beryllium	ND	0.00100	0.000176	1		Selenium	ND	0.0150	0.00295	1	
Cadmium	ND	0.00500	0.000350	1		Silver	ND	0.00500	0.000400	1	
Chromium	0.00650	0.00500	0.000350	1	В	Thallium	ND	0.0150	0.00233	1	
Cobalt	0.00194	0.00500	0.000696	1	J,B	Vanadium	0.00350	0.00500	0.000314	1 J	
Copper	ND	0.00500	0.00134	1		Zinc	0.0937	0.0100	0.000848	1	
Lead	ND	0.0100	0.00236	1							

MW-7 06-06-1241-3 06/20/06 Aqueous 06/21/06 06/22/06 060621L06F

Comment(s): -Results were evaluated to the MDL, concentrations >= to the MDL but < RL, if found, are qualified with a "J" flag.

-Mercury was analyzed on 6/22/2006 1:35:54 PM with batch 060621L04F

<u>Parameter</u>	<u>Result</u>	<u>RL</u>	MDL	<u>DF</u>	Qual	<u>Parameter</u>	Result	<u>RL</u>	<u>MDL</u>	DF Qual
Antimony	ND	0.0150	0.00209	1		Mercury	ND	0.000500	0.0000672	1
Arsenic	ND	0.0100	0.00308	1		Molybdenum	ND	0.00500	0.000800	1
Barium	0.146	0.010	0.000719	1		Nickel	0.00338	0.00500	0.00137	1 J,B
Beryllium	ND	0.00100	0.000176	1		Selenium	ND	0.0150	0.00295	1
Cadmium	ND	0.00500	0.000350	1		Silver	ND	0.00500	0.000400	1
Chromium	0.00641	0.00500	0.000350	1	В	Thallium	ND	0.0150	0.00233	1
Cobalt	0.00171	0.00500	0.000696	1	J,B	Vanadium	0.00393	0.00500	0.000314	1 J
Copper	ND	0.00500	0.00134	1		Zinc	0.110	0.010	0.000848	1
Lead	ND	0.0100	0.00236	1						

MW-3 06-06-1241-4 06/20/06 Aqueous 06/21/06; 06/22/06 060621L06F

-Results were evaluated to the MDL, concentrations >= to the MDL but < RL, if found, are qualified with a "J" flag.

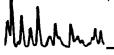
-Mercury was analyzed on 6/22/2006 1:38:05 PM with batch 060621L04F

D	Desuit	DI	A APOL	DE	Our	Daramatar	Desuit	Di	MOL	DE	O1
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	MDL	<u>DE</u>	<u>Qual</u>	<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>MDL</u>	DF 9	Quai
Antimony	ND	0.0150	0.00209	1		Mercury	ND	0.000500	0.0000672	1	
Arsenic	ND	0.0100	0.00308	1		Molybdenum	ND	0.00500	0.000800	1	
Barium	0.141	0.010	0.000719	1		Nickel	0.00365	0.00500	0.00137	1	J,B
Beryllium	ND	0.00100	0.000176	1		Selenium	ND	0.0150	0.00295	1	
Cadmium	ND	0.00500	0.000350	1		Silver	ND	0.00500	0.000400	1	
Chromium	0.00741	0.00500	0.000350	1	В	Thallium	0.00321	0.0150	0.00233	1	J
Cobalt	0.00197	0.00500	0.000696	1	J,B	Vanadium	0.00326	0.00500	0.000314	1	J
Copper	ND	0.00500	0.00134	1		Zinc	0.110	0.010	0.000848	1	
Lead	ND	0.0100	0.00236	1							

RL - Reporting Limit

DF - Dilution Factor

Qual - Qualifiers



Comment(s):





Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

06/20/06

Work Order No:

06-06-1241

Preparation:

EPA 3005A Filt. / EPA 7470A Filt.

Method:

EPA 6010B / EPA 7470A

Units:

mg/L

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Page 2 of 2

	Lab Sample	Date		Date	Date	
Client Sample Number	Number	Collected	Matrix	Prepared	Analyzed	QC Batch ID
Method Blank	099-04-008-2,530	NA NA	Aqueous	06/21/06	06/21/06	060621L04F

Comment(s): -Results were evaluated to the MDL, concentrations >= to the MDL but < RL, if found, are qualified with a "J" flag.

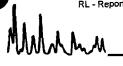
 Parameter
 Result
 RL

 Mercury
 ND
 0.000500

MDL D

DF Qual

Method Blank			097-	01-003-6,	222	N/A A	queous 06/21	/06 06/:	22/06 06062	1L06F	10 A
Comment(s):	-Results were eva	luated to the	MDL, concentrat	tions >= t	o the M	MDL but < RL, if four	nd, are qualified wit	h a "J" flag.			
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>MDL</u>	<u>DF</u>	<u>Qual</u>	<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>MDL</u>	DF Q	<u>ual</u>
Antimony	ND	0.0150	0.00209	1		Lead	ND	0.0100	0.00236	1	
Arsenic	ND	0.0100	0.00308	1		Molybdenum	ND	0.00500	0.000800	1	
Barium	ND	0.0100	0.000719	1		Nickel	0.00274	0.00500	0.00137	1	J
Beryllium	ND	0.00100	0.000176	1		Selenium	ND	0.0150	0.00295	1	
Cadmium	ND	0.00500	0.000350	1		Silver	0.000815	0.00500	0.000400	1	J
Chromium	0.00233	0.00500	0.000350	1	J	Thallium	ND	0.0150	0.00233	1	
Cobalt	0.00123	0.00500	0.000696	1	J	Vanadium	ND	0.00500	0.000314	1	
Copper	ND	0.00500	0.00134	1		Zinc	ND	0.0100	0.000848	1	







Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

Work Order No:

Preparation:

Method:

Units:

06/20/06 06-06-1241

EPA 3005A Filt.

EPA 6010B

mg/L

Project: BOLL Groundwater Monitoring

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602												
Client Sample No			Lab Nı	Sample Imber		Date Collected	Matrix		ate pared	Date Analyzed	Page 1 of	
MIVV-5			06-	06-1241-2		06/20/06	Aqueous	100	21/06	06/22/06	060621L06F	Courses A Service
Parameter Calcium Magnesium	Result 104 33.3	<u>RL</u> 0.100 0.1	MDL 0.00932 0.00328	DF 1 1 1 06-1241-3	Qual	Parameter Potassium Sodium 06/20/06		<u>sult</u> 5.74 37.7	RL 0.50 0.5	MDL 0.056 0.0192	<u>D</u> [Qual 1 B 1 B
_			<u> </u>			1.00 94 Co. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1.			1100	V6/22/06	060621L06F	
Parameter Calcium Magnesium	Result 98.3 34.4	RL 0.1 0.1	MDL 0.00932 0.00328	<u>DF</u> 1 1	Qual	<u>Parameter</u> Potassium Sodium		<u>sult</u> 5.45 7.9	<u>RL</u> 0.50 0.5	<u>MDL</u> 0.0561 0.0192	<u>DF</u> 1	Qual B B
MW-3		<u> </u>	06-0	6-1241-4		06/20/06	Aqueous	06/2	1/06	SAN KAMPUTTO VEZA	0 6 0621L06F	В
Parameter Calcium Magnesium	Result 107 32.7	<u>RL</u> 0.100 0.1	MDL 0.00932 0.00328	<u>DF</u> 9	Qual	Parameter Potassium Sodium	Res	<u>ult</u> 5.87 38.8	<u>RL</u> 0.50 0.5	MDL 0.0561 0.0192		В
Method Blank				01-003-6,2	Nach Addition	* NA	Aqueous	06/21	/06	06/22/06)60621L06F	В
Comment(s): Parameter Calcium Magnesium	-Results were eval <u>Result</u> ND ND	uated to the <u>RL</u> · 0.100 0.100	MDL, concentrate MDL 0.00932 0.00328	tions >≈ to <u>DF</u> (1 1	<u>auai</u>	DL but < RL, if fo <u>Parameter</u> Potassium Sodium	<u>Resi</u> 0.0		h a "J" fla RL 0.500 0.500	ag. <u>MDL</u> 0.0561 0.0192		· Qual J J

RL - Reporting Limit ,

DF - Dilution Factor

Qual - Qualifiers





Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

06/20/06

Work Order No:

06-06-1241

Preparation: Method:

EPA 3005A Filt. **EPA 200.8**

Units:

mg/L

Project: BO	U Groundwat		Pag	e 1 of 1							
Client Sample Nu	mber			ample nber		Date Collected	Matrix	Date Prepare		Date alyzed QC	Batch ID
MW-8			06-06	-1241-2		06/20/06	Aqueous	06/21/0)6 0	6/23/06 060	621L01F
Comment(s):	-Results were eva	luated to the	MDL, concentrate	tions >=	to the M	/IDL but < RL, if	found, are qua	lified with	a "J" flag		
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>MDL</u>	<u>DF</u>	<u>Qual</u>	<u>Parameter</u>	Res	<u>ult</u>	<u>RL</u>	<u>MDL</u>	DF Qual
Iron	0.0196	0.100	0.00214	1	J	Manganese	0.	00235	0.00100	0.0000189	1
MW-7			06-00	3-1241-3		06/20/06	Aqueous	06/21/0)6 (0	6/23/06 060	621L01F
Comment(s):	-Results were eva	luated to the	MDL, concentrate	ions >=	to the N	MDL but < RL, if	found, are qua	lified with	a "J" flag		
<u>Parameter</u>	Result	<u>RL</u>	MDL	<u>DF</u>	<u>Qual</u>	<u>Parameter</u>	<u>Res</u>	<u>ult</u>	<u>RL</u>	MDL	DF Qual
Iron	0.0210	0.100	0.00214	1	J	Manganese	0.	00184	0.00100	0.0000189	1
MW-3			06-01	5-1241-4		06/20/06	Aqueous	06/21/0)6 0	6/23/06 060	621L01F
Comment(s):	-Results were eva	luated to the	MDL, concentrate	tions >=	to the M	VIDL but < RL, if	found, are qua	lified with	a "J" flag		
<u>Parameter</u>	Result	<u>RL</u>	<u>MDL</u>	<u>DF</u>	<u>Qual</u>	<u>Parameter</u>	Res	<u>utt</u>	<u>RL</u>	<u>MDL</u>	DF Qual
Iron	0.0172	0.100	0.00214	1	J	Manganese	0.	00432	0.00100	0.0000189	_ 1
Method Blank			-099-	10-008-7	38	NA .	Aqueous	06/21/0)6 O	6/21/06 060	621L01F
Comment(s):	-Results were eva	luated to the	MDL, concentrate	tions >≃	to the 1	MDL but < RL, if	found, are qua	lified with	a "J" flag		
<u>Parameter</u>	Result	<u>RL</u>	MDL	<u>DF</u>	Qual	<u>Parameter</u>	<u>Res</u>	<u>ult</u>	<u>RL</u>	<u>MDL</u>	DF Qual
Iron	ND	0.100	0.00214	1		Manganese	N)	0.00100	0.0000189	1

RL - Reporting Limit ,

DF - Dilution Factor ,





Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

Work Order No:

Preparation:

Method:

06/20/06

06-06-1241

EPA 3520B EPA 8270C(M) Isotope

Dilution

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Page 1 of 1

Client Sample Number		Lab Samp Number		Date Collected	Matrix	Date Prepared	Date Analyzed	QC Batch ID
MW-8		06-06-12	41-2	06/20/06	Aqueous	06/22/06	06/28/06	060622L01D
Comment(s): -Results were	evaluated to the	MDL, concentrations	>= to the M	DL but < RL, if t	found, are qua	lified with a "J"	flag.	
<u>Parameter</u>	Result	<u>RL</u>	<u>MDL</u>	DF	<u>Qual</u>	<u>Units</u>		
1,4-Dioxane	ND	2.0	0.40	1		ug/L	٠.	
Surrogates:	<u>REC (%)</u>	Control Limits			<u>Qual</u>			
Nitrobenzene-d5	77	56-123						
MW-7	C. S. S.	06-06-12	41-3	06/20/06	Aqueous	06/22/06	06/28/06	∞ 060622L01D :
Comment(s): -Results were	evaluated to the	MDL, concentrations	>= to the M	DL but < RL, if	found, are qua	alified with a "J"	flag.	
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	MDL	<u>DF</u>	Qual	<u>Units</u>		
1,4-Dioxane	ND	2.0	0.40	1		ug/L		
Surrogates:	REC (%)	Control Limits			<u>Qual</u>			
Nitrobenzene-d5	85	56-123						
MW-3		06-06-12	41-4	06/20/06	Aqueous	06/22/06	06/28/06	060622L01D
Comment(s): -Results were	evaluated to the	MDL, concentrations	>≃ to the M	DL but < RL, if	found, are qua	alified with a "J"	flag.	
<u>Parameter</u>	Result	<u>RL</u>	<u>MDL</u>	DF	<u>Qual</u>	<u>Units</u>		
1,4-Dioxane	ND	2.0	0.40	1		ug/L		
Surrogates:	REC (%)	Control Limits			<u>Qual</u>			
Nitrobenzene-d5	81	56-123						
Method Blank		099-09-0	04-597	NA .	Aqueous	06/22/06	06/27/06	060622L01D
Comment(s): -Results were	evaluated to the						flag.	
<u>Parameter</u>	Result	<u>RL</u>	MDL	<u>DF</u>	Qual	<u>Units</u>		
	. 100	2.0	0.40	1		ug/L		
1,4-Dioxane	ND		0.70	-		3		
1,4-Dioxane Surrogates:	REC (%)	Control Limits	0.10	•	Qual	-		

DF - Dilution Factor ,

Qual - Qualifiers





Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

Work Order No:

Preparation:

Method:

06/20/06

06-06-1241

EPA 3520B

EPA 1625CM

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Page 1 of 1

Project: BOU Ground	dwater Mon	itoring 2006 (F	AC Wel	ls) / 1/653	-0602			Page 1 of 1
Client Sample Number		Lab Sam Numbe		Date Collected	Matrix	Date Prepared	Date Analyzed	QC Batch ID
MW-8	*	06-06-12	241-2	06/20/06	Aqueous	06/23/06	06/27/06	060623L09
Comment(s): -Results were	evaluated to the	MDL, concentrations	s >= to the M	IDL but < RL, if	found, are qua	alified with a "J"	flag.	
Parameter	Result	RL	MDL	<u>DF</u>	<u>Qual</u>	<u>Units</u>		
N-Nitrosodimethylamine	ND	2.0	0.48	1 1		ng/L		
urrogates:	<u>REC (%)</u>	Control Limits			<u>Qual</u>			
,4-Dichlorobenzene-d4	53	50-130						
MW-7:	18	06-06-12	241-3	06/20/06	Aqueous	06/23/06	06/27/06	- 060623L09
Comment(s): -Results were		•					flag.	
arameter	Result	<u>RL</u>	MDL	<u>DF</u>	<u>Qual</u>	<u>Units</u>		•
l-Nitrosodimethylamine	ND	2.0	0.48	1		ng/L		
urrogates:	<u>REC (%)</u>	Control Limits			<u>Qual</u>			
,4-Dichlorobenzene-d4	52	50-130						
MW-3		06-06-12	241-4	06/20/06	Aqueous	06/23/06	06/27/06	060623L09
Comment(s): -Results were	evaluated to the		s >= to the M	•	found, are qua		flag.	
arameter	Result	<u>RL</u>	MDL	<u>DF</u>	<u>Qual</u>	<u>Units</u>		
I-Nitrosodimethylamine	ND	2.0	0.48	1		ng/L		
urrogates:	REC (%)	Control Limits			Qual			
,4-Dichlorobenzene-d4	50	50-130						
Method Blank		3.5 j 099-07-0	27-251	NA . A	#-Aqueous	06/23/06	06/27/06	- 060623L09
Comment(s): -Results were		•					flag.	
<u>'arameter</u>	Result	<u>RL</u>	<u>MDL</u>	<u>DF</u>	Qual	<u>Units</u>		
I-Nitrosodimethylamine	ND	2.0	0.48	1		ng/L		
urrogates:	<u>REC (%)</u>	Control Limits			<u>Qual</u>			
,4-Dichlorobenzene-d4	54	50-130						

RL - Reporting Limit

DF - Dilution Facto

Qual - Qualifier





Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

Work Order No:

Preparation:

Method:

Units:

06/20/06 06-06-1241

EPA 5030B

EPA 8260B

ug/L

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Page 1 of 6

Client Sample Number		•	Lab Sa Numl			Date Collected Matrix	Date Prepared	Date Analyzed	QC Bate	th ID	-
LTB062006		186	06-06-	1241-1		06/20/06 Aqueous	07/03/06	07/03/06	060703	L 01	
Comment(s): -Results were	evaluated to th	e MDL, co	ncentratio	ns >= to	the N	ADL but < RL, if found, are quality	fied with a "J"	flag.			
<u>Parameter</u>	Result	<u>RL</u>	<u>MDL</u>	DF Q	ual	<u>Parameter</u>	Result	RL	<u>MDL</u>	<u>D</u> F	Qual
Acetone	9.4	10.0	6.1	1,	J,B	1,3-Dichloropropane	ND	1.0	0.30	1	
Benzene	ND	0.50	0.26	1		2,2-Dichloropropane	ND	1.0	0.40	1	
Bromobenzene	ND	1.0	0.47	1		1,1-Dichloropropene	ND	1.0	0.21	1	
Bromochloromethane	МD	1.0	0.68	1		c-1,3-Dichloropropene	ND	0.50	0.45	1	
Bromodichloromethane	ND	1.0	0.27	1		t-1,3-Dichloropropene	ND	0.50	0.31	1	
Bromoform	ND	1.0	0.62	1		Ethylbenzene	ND	1.0	0.17	1	
Bromomethane	ND	10	2.9	1		2-Hexanone	ND	10	1.9	1	
2-Butanone	ND	10	4.2	1		Isopropylbenzene	ND	1.0	0.24	1	
n-Butylbenzene	ND	1.0	0.29	1		p-Isopropyttoluene	ND	1.0	0.21	1	
sec-Butylbenzene	ND	1.0	0.21	1		Methylene Chloride	3.4	10.0	2.6	1	J,B
tert-Butylbenzene	ND	1.0	0.17	1		4-Methyl-2-Pentanone	ND	10	2.4	1	
Carbon Disulfide	ND	10	1.0	1		Naphthalene	ND	10	0.95	1	
Carbon Tetrachloride	ND	0.50	0.42	1		n-Propylbenzene	ND	1.0	0.30	1	(
Chlorobenzene	ND	1.0	0.36	1		Styrene	ND	1.0	0.29	1	
Chloroethane	ND	1.0	0.52	1		1,1,1,2-Tetrachloroethane	ND	1.0	0.37	1	
Chloroform	ND	1.0	0.22	1		1,1,2,2-Tetrachloroethane	ND	1.0	0.37	1	
Chloromethane	ND	10	1.8	1		Tetrachloroethene	ND	1.0	0.29	1	
2-Chlorotoluene	ND	1.0	0.24	1		Toluene	0.41	1.0	0.35	1	J
4-Chlorotoluene	ND	1.0	0.30	1		1,2,3-Trichlorobenzene	ND	1.0	0.39	1	
Dibromochloromethane	ND	1.0	0.45	1		1,2,4-Trichlorobenzene	ND	1.0	0.35	1	
1.2-Dibromo-3-Chloropropane	ND	5.0	2.5	1		1,1,1-Trichloroethane	ND	1.0	0.32	1	
1,2-Dibromoethane	ND	1.0	0.81	1		1,1,2-Trichloro-1,2,2-Trifluoroe	thane ND	10	0.54	1	
Dibromomethane	ND	1.0	0.42	1		1,1,2-Trichloroethane	ND	1.0	0.54	1	
1,2-Dichlorobenzene	ПD	1.0	0.24	1		Trichloroethene	ND	1.0	0.30	1	
1,3-Dichlorobenzene	ND	1.0	0.38	1		Trichlorofluoromethane	ND	10	0.36	1	
1,4-Dichlorobenzene	ND	1.0	0.30	1		1,2,3-Trichloropropane	ND	5.0	2.3	1	
Dichlorodifluoromethane	ND	1.0	0.27	1		1,2,4-Trimethylbenzene	ND	1.0	0.26	1	
1,1-Dichloroethane	ND	1.0	0.53	1 ,		1,3,5-Trimethylbenzene	ND	1.0	0.19	1	
1,2-Dichloroethane	ND	0.50	0.22	1		Vinyl Acetate	ND	10	3.2	1	
1,1-Dichloroethene	ND	1.0	0.31	1		Vinyl Chloride	ND	0.50	0.33	1	
c-1,2-Dichloroethene	ND	1.0	0.35	1		p/m-Xylene	ND	1.0	0.38	1	
t-1,2-Dichloroethene	ND	1.0	0.29	1		o-Xylene	ND	1.0	0.21	1	
1,2-Dichloropropane	ND	1.0	0.28	1		Methyl-t-Butyl Ether (MTBE)	ND	1.0	0.29	1	
Surrogates:	<u>REC (%)</u>	Control	Limits ·	<u>Q</u>	ual	Surrogates:	<u>REC (%</u>	6) Control	Limits		Qual
Dibromofluoromethane	108	74-140				1,2-Dichloroethane-d4	101	74-146			
Toluene-d8	97	88-112				1,4-Bromofluorobenzene	89	74-110			







Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

06/20/06

Work Order No:

Date

Date

06-06-1241

Preparation:

EPA 5030B

Method:

Date

EPA 8260B

Units:

ug/L

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Lab Sample

Page 2 of 6

Client Sample Number			Numt	er		Collected Matrix P	repared	Analyzed	QC Bat	ch ID
MW-8			06-06-	1241-2		06/20/06 Aqueous 0	7/03/06	07/03/06	060703	L01
Comment(s): -Results were	evaluated to th	ne MDL, co	ncentratio	ns >= to	the N	MDL but < RL, if found, are qualified	d with a "J" f	lag.		
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	MDL	DF 9	Qual	Parameter	<u>Result</u>	<u>RL</u>	MDL	DF Qual
Acetone	10	10	6.1	1	В	1,3-Dichloropropane	ND	1.0	0.30	1
Benzene	ND	0.50	0.26	1		2,2-Dichloropropane	ND	1.0	0.40	1
Bromobenzene	ND	1.0	0.47	1		1,1-Dichloropropene	ND	1.0	0.21	1
Bromochloromethane	ND	1.0	0.68	1		c-1,3-Dichloropropene	ND	0.50	0.45	1
Bromodichloromethane	ND	1.0	0.27	1		t-1,3-Dichloropropene	ND	0.50	0.31	1
Bromoform	ND	1.0	0.62	1		Ethylbenzene	ND	1.0	0.17	1
Bromomethane	ND	10	2.9	1		2-Hexanone	ND	10	1.9	1
2-Butanone	ND	10	4.2	1		Isopropylbenzene	ND	1.0	0.24	1 -
n-Butylbenzene	ND	1.0	0.29	1		p-Isopropyltoluene	ND	1.0	0.21	1
sec-Butylbenzene	ND	1.0	0.21	1		Methylene Chloride	4.2	10.0	2.6	1 J,B
tert-Butylbenzene	ND	1.0	0.17	1		4-Methyl-2-Pentanone	ND	10	2.4	1
Carbon Disulfide	ND	10	1.0	1		Naphthalene	ND	10	0.95	1
Carbon Tetrachloride	0.79	0.50	0.42	1		n-Propylbenzene	ND	1.0	0.30	1
Chlorobenzene	ND	1.0	0.36	1		Styrene	ND	1.0	0.29	1
Chloroethane	ND	1.0	0.52	1		1,1,1,2-Tetrachloroethane	ND	1.0	0.37	1
Chloroform	1.1	1.0	0.22	1		1,1,2,2-Tetrachloroethane	ND	1.0	0.37	1
Chloromethane	ND	10	1.8	1		Tetrachloroethene	160	1	0.29	1 .
2-Chlorotoluene	ND	1.0	0.24	1		Toluene	ND	1.0	0.35	1
4-Chlorotoluene	ND	1.0	0.30	1		1,2,3-Trichlorobenzene	ND	1.0	0.39	1
Dibromochloromethane	ND	1.0	0.45	1		1,2,4-Trichlorobenzene	ND	1.0	0.35	1
1,2-Dibromo-3-Chloropropane	ND	5.0	2.5	1		1,1,1-Trichloroethane	ND	1.0	0.32	1
1,2-Dibromoethane	ND	1.0	0.81	1		1,1,2-Trichloro-1,2,2-Trifluoroeth		10.0	0.54	1 J
Dibromomethane	ND	1.0	0.42	1		1,1,2-Trichloroethane	ND	1.0	0.54	1
1,2-Dichlorobenzene	ND	1.0	0.24	1		Trichloroethene	58	1	0.30	1
1,3-Dichlorobenzene	ND	1.0	0.38	1		Trichlorofluoromethane	ND	10	0.36	1
1,4-Dichlorobenzene	ND	1.0	0.30	1		1,2,3-Trichloropropane	ND	5.0	2.3	1
Dichlorodifluoromethane	ND	1.0	0.27	1		1,2,4-Trimethylbenzene	ND	1.0	0.26	1
1,1-Dichloroethane	ND	1.0	0.53	1		1,3,5-Trimethylbenzene	ND	1.0	0.19	1
1,2-Dichloroethane	1.6	0.5	0.22	1		Vinyl Acetate	ND	10	3.2	1
1,1-Dichloroethene	0.97	1.0	0.31	1	J	Vinyl Chloride	ND	0.50	0.33	1
c-1,2-Dichloroethene	ND	1.0	0.35	1		p/m-Xylene	ND	1.0	0.38	1
t-1,2-Dichloroethene	ND	1.0	0.29	1		o-Xylene	ND	1.0	0.21	1
1,2-Dichloropropane	ND	1.0	0.28	1		Methyl-t-Butyl Ether (MTBE)	ND	1.0	0.29	1







Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

Work Order No:

Preparation:

Method:

Date

Units:

06/20/06 06-06-1241 **EPA 5030B**

EPA 8260B

ug/L

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

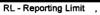
Lab Sample

Page 3 of 6

Date

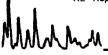
Date

Client Sample Number		•	Lab Sai Numb		Collected Matrix	Prepared	Date Analyzed	QC Bate	ch ID
MW-7			06-06-	1241-3	: 06/20/06 Aqueous	07/03/06	07/03/06	060703	L01
Comment(s): -Results were e	valuated to the	MDL, co	ncentratio	ns >= to the	MDL but < RL, if found, are qual	ified with a "J" fl	ag.		
<u>Parameter</u>	Result	<u>RL</u>	MDL	DF Qual	· · · · · · · · · · · · · · · · · · ·	Result	<u>RL</u>	MDL	DF Qual
Acetone	9.8	10.0	6.1	1 J,B	1,3-Dichloropropane	ND	1.0	0.30	1
Benzene	ND	0.50	0.26	1	2,2-Dichloropropane	ND	1.0	0.40	1
Bromobenzene	ND	1.0	0.47	1	1.1-Dichloropropene	ND	1.0	0.21	1
Bromochloromethane	ND	1.0	0.68	1	c-1,3-Dichloropropene	ND	0.50	0.45	1
Bromodichloromethane	ND	1.0	0.27	1	t-1,3-Dichloropropene	ND	0.50	0.31	1
Bromoform	ND	1.0	0.62	1	Ethylbenzene	ND	1.0	0.17	1
Bromomethane	ND	10	2.9	1	2-Hexanone	NĐ	10	1.9	1
2-Butanone	ND	10	4.2	1	Isopropylbenzene	ND	1.0	0.24	1
n-Butylbenzene	ND	1.0	0.29	1	p-Isopropyltoluene	ND	1.0	0.21	1
sec-Butylbenzene	ND	1.0	0.21	1	Methylene Chloride	3.6	10.0	2.6	1 J,B
tert-Butylbenzene	ND	1.0	0.17	1	4-Methyl-2-Pentanone	ND	10	2.4	1
Carbon Disulfide	ND	10	1.0	1	Naphthalene	ND	10	0.95	1 🚄
Carbon Tetrachloride	ND	0.50	0.42	1	n-Propylbenzene	ND	1.0	0.30	1
Chlorobenzene	ND	1.0	0.36	1	Styrene	ND	1.0	0.29	1
Chloroethane	ND	1.0	0.52	1	1,1,1,2-Tetrachloroethane	ND	1.0	0.37	1
Chloroform	0.71	1.0	0.22	1 · J	1,1,2,2-Tetrachloroethane	ND	1.0	0.37	1
Chloromethane	ND	10	1.8	1	Tetrachloroethene	50	1	0.29	1
2-Chiorotoluene	ND	1.0	0.24	1	Toluene	ND	1.0	0.35	1
4-Chlorotoluene	ND	1.0	0.30	1	1,2,3-Trichlorobenzene	ND	1.0	0.39	1
Dibromochloromethane	ND	1.0	0.45	1	1,2,4-Trichlorobenzene	ND	1.0	0.35	1
1,2-Dibromo-3-Chloropropane	ND	5.0	2.5	1	1,1,1-Trichloroethane	ND	1.0	0.32	1
1,2-Dibromoethane	ND	1.0	0.81	1	1,1,2-Trichloro-1,2,2-Trifluoro	ethane ND	10	0.54	1
Dibromomethane	ND	1.0	0.42	1	1,1,2-Trichloroethane	ND	1.0	0.54	1
1,2-Dichlorobenzene	ND	1.0	0.24	1	Trichloroethene	17	1	0.30	1
1,3-Dichlorobenzene	ND	1.0	0.38	1	Trichlorofluoromethane	ND	10	0.36	1
1,4-Dichlorobenzene	ND	1.0	0.30	1	1,2,3-Trichloropropane	ND	5.0	2.3	1
Dichlorodifluoromethane	ND	1.0	0.27	1	1,2,4-Trimethylbenzene	ND	1.0	0.26	1
1,1-Dichloroethane	ND	1.0	0.53	1	1,3,5-Trimethylbenzene	ND	1.0	0.19	1
1,2-Dichloroethane	ND	0.50	0.22	1	Vinyl Acetate	ND	10	3.2	1
1,1-Dichloroethene	0.41	1.0	0.31	1 J	Vinyl Chloride	, ND	0.50	0.33	1
c-1,2-Dichloroethene	ND	1.0	0.35	1	p/m-Xylene	ND	1.0	0.38	1
t-1,2-Dichloroethene	ND	1.0	0.29	1	o-Xylene	ND	1.0	0.21	1
1,2-Dichloropropane	ND	1.0	0.28	1	Methyl-t-Butyl Ether (MTBE)	ND	1.0	0.29	1
Surrogates:	REC (%)	Control	<u>Limits</u>	<u>Qual</u>	Surrogates:	<u>REC (%</u>) Control	<u>Limits</u>	<u>Qual</u>
Dibromofluoromethane	109	74-140			1,2-Dichloroethane-d4	102	74-146		
Toluene-d8	97	88-112			1,4-Bromofluorobenzene	87	74-110		



DF - Dilution Factor

Qual - Qualifiers







Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

Work Order No:

Date

Date

Preparation:

Method:

Units:

Date

06-06-1241 EPA 5030B

06/20/06

EPA 8260B

ug/L

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Lab Sample

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Client Sample Number			Lab Sai Numb			Date Collected	Matrix	Date Prepared	Date Analyzed	QC Bate	ch ID
MW-3		3	, 06-06-	1241-4		06/20/06%	Aqueous	07/03/06	07/03/06	060703	L01
Comment(s): -Results were en	valuated to th	e MDL, co	ncentratio	ons >= t	to the N	MDL but < RL, if f	ound, are quali	fied with a "J" i	flag.		
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	MDL	<u>DF</u>	<u>Qual</u>	<u>Parameter</u>		Result	<u>RL</u>	MDL.	DF Qual
Acetone	9.7	10.0	6.1	1	J,B	1,3-Dichloropro	pane	ND	1.0	0.30	1
Benzene	ND	0.50	0.26	1 1		2,2-Dichloropro		ND	1.0	0.40	1
Bromobenzene	ND	1.0	0.47	1		1,1-Dichloropro	pene	ND	1.0	0.21	1
Bromochloromethane	ND	1.0	0.68	1		c-1,3-Dichlorop	ropene	ND	0.50	0.45	1 .
Bromodichloromethane	ND	1.0	0.27	1		t-1,3-Dichloropr		ND	0.50	0.31	1
Bromoform	ND	1.0	0.62	1		Ethylbenzene	•	ND	1.0	0.17	1
Bromomethane	ND	10	2.9	1		2-Hexanone		ND	10	1.9	1
2-Butanone	ND	10	4.2	1		isopropylbenze	ne	ND	1.0	0.24	1
n-Butylbenzene	ND	1.0	0.29	1		p-isopropyttolue	ene	ND	1.0	0.21	1
sec-Butylbenzene	ND	1.0	0.21	1		Methylene Chlo	ride	3.6	10.0	2.6	1 J,B
tert-Butylbenzene	ND	1.0	0.17	1		4-Methyl-2-Pen	tanone	ND	10	2.4	1
Carbon Disulfide	ND	10	1.0	1		Naphthaiene		ND	10	0.95	1
Carbon Tetrachloride	1.1	0.5	0.42	1		n-Propylbenzen	e	ND	1.0	0.30	1
Chlorobenzene	ND	1.0	0.36	1		Styrene		ND	1.0	0.29	1
Chloroethane	ND	1.0	0.52	1		1,1,1,2-Tetrach	loroethane	ND	1.0	0.37	1.1
Chloroform	1.4	1.0	0.22	1		1,1,2,2-Tetrach	loroethane	ND	1.0	0.37	1
Chloromethane	ND	10	1.8	1		Tetrachloroethe	ene	36	1	0.29	1
2-Chlorotoluene	ND	1.0	0.24	1		Toluene		ND	1.0	0.35	1
4-Chlorotoluene	ND	1.0	0.30	1		1,2,3-Trichlorob	penzene	ND	1.0	0.39	1
Dibromochloromethane	ND	1.0	0.45	1		1,2,4-Trichlorob	penzene	ND	1.0	0.35	1
1,2-Dibromo-3-Chloropropane	ND	5.0	2.5	1		1,1,1-Trichloroe	ethane	ND	1.0	0.32	1
1,2-Dibromoethane	ND	1.0	0.81	1		1,1,2-Trichloro-	1,2,2-Trifluoroe	ethane 1.9	10.0	0.54	1 J
Dibromomethane	ND	1.0	0.42	1		1,1,2-Trichloroe	ethane	ND .	1.0	0.54	1
1,2-Dichtorobenzene	ND	1.0	0.24	1		Trichloroethene	•	12	1	0.30	1
1,3-Dichlorobenzene	ND	1.0	0.38	1		Trichlorofluoron	nethane	ND	10	0.36	1
1,4-Dichlorobenzene	ND	1.0	0.30	1		1,2,3-Trichlorop		ND	5.0	2.3	1
Dichlorodifluoromethane	ND	1.0	0.27	1		1,2,4-Trimethyll		ND	1.0	0.26	1
1,1-Dichloroethane	ND	1.0	0.53	1		1,3,5-Trimethyll	benzene	ND	1.0	0.19	1
1,2-Dichloroethane	0.46	0.50	0.22	1	J	Vinyl Acetate		ND	10	3.2	1
1,1-Dichloroethene	2.6	1.0	0.31	1		Vinyl Chloride		ND	0.50	0.33	1
c-1,2-Dichloroethene	ND	1.0	0.35	1		p/m-Xylene		ND	1.0	0.38	1
t-1,2-Dichloroethene	ND	1.0	0.29	1		o-Xylene		ND	1.0	0.21	1
1,2-Dichloropropane	ND .	1.0	0.28	1		Methyl-t-Butyl E	ther (MTBE)	ND O	1.0	0.29	1
Surrogates:	<u>REC (%)</u>	Control I	<u>imits</u>		<u>Qual</u>	Surrogates:	•	<u>REC (%</u>		<u>Limits</u>	<u>Qual</u>
Dibromofluoromethane	110	74-140				1,2-Dichloroeth		103	74-146		
Toluene-d8	98	88-112				1,4-Bromofluoro	obenzene	86	74-110		







Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300

3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Date Received:

06/20/06

Work Order No: Preparation:

Date

Date

06-06-1241 EPA 5030B

Method:

EPA 8260B

Units:

Date

ug/L

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Lab Sample

Page 5 of 6

Client Sample Number			Numb	ег	Collected	Matrix	Prepared	Analyzed	QC Bate	th ID	
LFB-062006			06-06-	1241-5	06/20/06	Aqueous	07/03/06	07/03/06	060703	L01 ·	
Comment(s): -Results were	evaluated to the		ncentratio		MDL but < RL, if	found, are quali	fied with a "J" f	-			
<u>Parameter</u>	Result	<u>RL</u>	<u>MDL</u>	DF Qua	<u>Parameter</u>		<u>Result</u>	<u>RL</u>	<u>MDL</u>	<u>DF</u> Q	≀ual
Acetone	9.9	10.0	6.1	1 J,I	1,3-Dichloropr	opane	ND	1.0	0.30	1	
Benzene	ND	0.50	0.26	1	2,2-Dichloropr	opane	ND	1.0	0.40	1	
Bromobenzene	ND	1.0	0.47	1	1,1-Dichloropr	opene	ND	1.0	0.21	1	
Bromochloromethane	ND	1.0	0.68	1	c-1,3-Dichloro	propene	ND	0.50	0.45	1	
Bromodichtoromethane	ND	1.0	0.27	1	t-1,3-Dichlorop	oropene	ND	0.50	0.31	1	
Bromoform	ND	1.0	0.62	1	Ethylbenzene		ND	1.0	0.17	1	
Bromomethane	ND	10	2.9	1	2-Hexanone		ND	10	1.9	1	
2-Butanone	ND	10	4.2	1	Isopropylbenze	ene	ND	1.0	0.24	1	
n-Butylbenzene	ND	1.0	0.29	1	p-Isopropyitoli	Jene	ND	1.0	0.21	1	•
sec-Butylbenzene	ND	1.0	0.21	1	Methylene Chi	loride	5.5	10.0	2.6	1	J,B
tert-Butylbenzene	ND	1.0	0.17	1	4-Methyl-2-Pe	ntanone	ND	10	2.4	1	
Carbon Disulfide	ND	10	1.0	1	Naphthalene		ND	10	0.95	1	- 1
Carbon Tetrachloride	ND	0.50	0.42	1	n-Propylbenze	ene	ND	1.0	0.30	1	
Chlorobenzene	ND	1.0	0.36	1	Styrene		ND	1.0	0.29	1	
Chloroethane	ND	1.0	0.52	1	1,1,1,2-Tetrac	hloroethane	ND	1.0	0.37	1	
Chloroform	ND	1.0	0.22	1	1,1,2,2-Tetrac	hioroethane	ND	1.0	0.37	1	
Chloromethane	ND	10	1.8	1	Tetrachloroeth	nene	ND	1.0	0.29	1	
2-Chlorotoluene	ND	1.0	0.24	1	Toluene		ND	1.0	0.35	1	
4-Chlorotoluene	ND	1.0	0.30	1	1,2,3-Trichlore	benzene	ND	1.0	0.39	1	
Dibromochloromethane	ND	1.0	0.45	1	1,2,4-Trichlore	benzene	ND	1.0	0.35	1	
1,2-Dibromo-3-Chloropropane	ND	5.0	2.5	. 1	1,1,1-Trichlore	ethane	ND	1.0	0.32	1	
1,2-Dibromoethane	ND	1.0	0.81	1	1,1,2-Trichlore	-1,2,2-Trifluoro	ethane ND	10	0.54	1	
Dibromomethane	ND	1.0	0.42	1	1,1,2-Trichlore	ethane	ND	1.0	0.54	1	
1,2-Dichlorobenzene	ND	1.0	0.24	1	Trichloroether	ne	ND	1.0	0.30	1	
1,3-Dichlorobenzene	ND	1.0	0.38	1	Trichlorofluoro	methane	ND	10	0.36	1	
1,4-Dichlorobenzene	ND	1.0	0.30	1	1,2,3-Trichlore	opropane	ND	5.0	2.3	1	
Dichlorodifluoromethane	ND	1.0	0.27	1	1,2,4-Trimethy		ND	1.0	0.26	1	
1,1-Dichloroethane	ND	1.0	0.53	1	1,3,5-Trimethy	/ibenzene	ND	1.0	0.19	1	
1,2-Dichloroethane	ND	0.50	0.22	1	Vinyl Acetate		ND	10	3.2	1	
1,1-Dichloroethene	ND	1.0	0.31	1	Vinyl Chloride		·ND	0.50	0.33	1	
c-1,2-Dichloroethene	ND	1.0	0.35	1	p/m-Xylene		ND	1.0	0.38	1	
t-1,2-Dichtoroethene	ND	1.0	0.29	1	o-Xylene		ND	1.0	0.21	1	
1,2-Dichloropropane	ND	1.0	0.28	1		Ether (MTBE)	ND	1.0	0.29	1	
Surrogates:	<u>REC (%)</u>	Control	<u>Limits</u>	. Qu	Surrogates:		<u>REC (%</u>	<u>(6)</u> <u>Control</u>	<u>Limits</u>	Q	Qual
Dibromofluoromethane	110 ·	74-140			1,2-Dichloroet	hane-d4	106	74-146			
Toluene-d8	99	88-112			1,4-Bromofluo	robenzene	88	74-110			



DF - Dilution Factor

Qual - Qualifiers







Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received:

Work Order No:

Date

Date

Preparation: Method:

Units:

Date

06-06-1241 **EPA 5030B**

06/20/06

EPA 8260B

ug/L

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Lab Sample

Page 6 of 6

Client Sample Number			Num		Collected	Matrix	Prepared	Analyzed	QC Bat	ch ID
		100	099_1	0-006-18,441/	N/A	Aqueous	07/03/06	.07/03/06	060703	101
Metrou Dialik						Adresona	01/05/00	,0,,00,00		LU
Comment(s): -Results were ev	aluated to the	e MDL, coi	ncentratio	ons >= to the N	MDL but < RL, if f	found, are qualifi	ed with a "J" f	lag.		
Parameter	<u>Result</u>	RL	MDL	DF Qual	Parameter	•	Result	RL	MDL	DF Qual
Acetone	12	10	6.1	1	1,3-Dichtoropro	pane	ND	1.0	0.30	1
Benzene	ND	0.50	0.26	1	2,2-Dichloropro		ND	1.0	0.40	1
Bromobenzene	ND	1.0	0.47	1	1,1-Dichloropro		ND	1.0	0.21	1
Bromochloromethane	ND	1.0	0.68	1	c-1,3-Dichlorop	•	ND	0.50	0.45	1
Bromodichloromethane	ND	1.0	0.27	1	t-1,3-Dichlorop	•	ND	0.50	0.31	1
Bromoform	ND	1.0	0.62	1	Ethylbenzene	•	ND	1.0	0.17	1
Bromomethane	ND	10	2.9	1	2-Hexanone		ND	10	1.9	1
2-Butanone	ND	10	4.2	1	Isopropylbenze	ene	ND	1.0	0.24	1
n-Butylbenzene	ND	1.0	0.29	1	p-Isopropyltolu	ene	ND	1.0	0.21	1
sec-Butylbenzene	ND	1.0	0.21	1	Methylene Chlo	oride	3.2	10.0	2.6	1 J
tert-Butylbenzene	ND	1.0	0.17	1	4-Methyl-2-Per	ntanone	ND	10	2.4	1
Carbon Disulfide	ND	10	1.0	1	Naphthalene		ND	10	0.95	1
Carbon Tetrachloride	ND	0.50	0.42	1	n-Propylbenzer	ne	ND	1.0	0.30	1
Chlorobenzene	ND	1.0	0.36	1	Styrene		ND	1.0	0.29	1
Chloroethane	ND	1.0	0.52	1	1,1,1,2-Tetrach	nloroethane	ND	1.0	0.37	1
Chloroform	ND	1.0	0.22	1	1,1,2,2-Tetrach	nloroethane	ND	1.0	0.37	1
Chloromethane	ND	10	1.8	1	Tetrachloroethe	ene	ND	1.0	0.29	1
2-Chlorotoluene	ND	1.0	0.24	1	Toluene		ND	1.0	0.35	1
4-Chlorotoluene	ND	1.0	0.30	1	1,2,3-Trichlorol	benzene	ND	1.0	0.39	1
Dibromochloromethane	ND	1.0	0.45	1	1,2,4-Trichlorol	benzene	ND	1.0	0.35	1
1,2-Dibromo-3-Chloropropane	ND	5.0	2.5	1	1,1,1-Trichloro	ethane	ND	1.0	0.32	1
1,2-Dibromoethane	ND	1.0	0.81	1	1,1,2-Trichloro-	-1,2,2-Trifluoroet	thane ND	10	0.54	1
Dibromomethane	ND	1.0	0.42	1	1,1,2-Trichloro	ethane	ND	1.0	0.54	1
1,2-Dichlorobenzene	ND	1.0	0.24	1	Trichloroethene	е	ND	1.0	0.30	1
1,3-Dichlorobenzene	ND	1.0	0.38	1	Trichlorofluoror	methane	ND	10	0.36	1
1,4-Dichlorobenzene	ND	1.0	0.30	1	1,2,3-Trichloro	propane	ND	5.0	2.3	1
Dichlorodifluoromethane	ND	1.0	0.27	1	1,2,4-Trimethyl		ND	1.0	0.26	1
1,1-Dichloroethane	ND	1.0	0.53	1	1,3,5-Trimethyl	ibenzene	ND	1.0	0.19	1
1,2-Dichloroethane	ND	0.50	0.22	1	Vinyl Acetate		ND	10	3.2	1
1,1-Dichloroethene	ND	1.0	0.31	1	Vinyl Chloride		ND	0.50	0.33	1
c-1,2-Dichloroethene	ND	1.0	0.35	1	p/m-Xylene		ND	1.0	0.38	1
t-1,2-Dichloroethene	ND	1.0	0.29	1	o-Xylene		ND	1.0	0.21	1
1,2-Dichloropropane	ND	1.0	0.28	1	Methyl-t-Butyl E	Ether (MTBE)	ND ND	1.0	0.29	1
Surrogates:	<u>REC (%)</u>	Control L	<u>limits</u>	<u>Qual</u>	Surrogates:		<u>REC (%</u>		Limits	<u>Quat</u>
Dibromofluoromethane	105	74-140			1,2-Dichloroeth		98	74-146		
Toluene-d8	97	88-112			1,4-Bromofluor	robenzene	89	74-110		



DF - Dilution Factor ,



EPA 8260B Tentatively Identified Compound List

Work Order CEL Sample Client ID

Q Compound

CAS NUMBER

<u>RT</u>

On Column Conc. Estimated Conc.

ug/L

ug/L

06-06-1241

No TiCs found for all samples

Q Qualifier RT Retention Time

Mulum

Page 1 of 1





Tetra Tech, Inc. 3475 East Foothill Blvd., Suite

3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Date Received: Work Order No: Preparation:

Method:

06/20/06 06-06-1241 EPA 5030B SRL 524M-TCP

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Page 1 of 1

Client Sample Number		Lab Sar Numb		Date Collected	Matrix	Date Prepared	Date Analyzed	QC Batch ID
MW-8	ar e	06-06-	1241-2	06/20/06	Aqueous	906/21/06	06/22/06	060621L02
<u>Parameter</u>	Result	<u>RL</u>	MDL	<u>DF</u>	<u>Qual</u>	<u>Units</u>		
1,2,3-Trichloropropane	0.068	0.005	0.0017	1		ug/L		
MW-7		06-06-	1241-3	06/20/06	Aqueous	06/21/06	06/22/06	0606211.02
<u>Parameter</u>	Result	RL	MDL	<u>DF</u>	Qual	<u>Units</u>		
1,2,3-Trichloropropane	0.016	0.005	0.0017	1		ug/L		
MW-3		06-06-	1241-4	06/20/06	Aqueous	[®] 06/21/06	06/22/06	060621L02
<u>Parameter</u>	Result	<u>RL</u>	MDL	<u>DF</u>	Quat	<u>Units</u>		
1,2,3-Trichloropropane	0.044	0.005	0.0017	1		ug/L		
Method Blank		099-10	-022-240	. NA	Aqueous	06/21/06	06/22/06	060621L02
Comment(s): -Results were	evaluated to the I	MDL, concentratio	ns >= to the Mi	OL but < RL, if	found, are qua	alified with a "J"	flag.	
<u>Parameter</u>	Result	RL	<u>MDL</u>	<u>DF</u>	Qual	<u>Units</u>		
1,2,3-Trichloropropane	ND	0.0050	0.0017	1		ug/L		

Lh.u_

DF - Dilution Factor ,

Qual - Qualifiers





Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received:

Work Order No:

06/20/06

06-06-1241

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Page 1 of 2

		L	ab Sample	Number		Matri	i.,		
Client Sample Num	oer				Collected	IVIAL	IX	·	
MW-8			06-06-124	1-2	06/20/06	Aque	ous	1000	
Comment(s): (1)	Results were evaluated to the	e MDL, conce	entrations >	= to the	MDL but <	RL, if found	i, are qualified with	n a "J" flag.	
<u>Parameter</u>	Result	<u>RL</u>	MDL.	<u>DF</u>	Qual	<u>Units</u>	Date Prepared	Date Analyzed	<u>Method</u>
Chromium, Hexaval	ent 1.5	0.2	0.0050	1	В	ug/L	N/A	06/20/06	EPA 218.6
Chloride	40	10	0.055	10		mg/L	N/A	06/21/06	EPA 300.0
Nitrite (as N) (1)	ND	0.10	0.015	1		mg/L	N/A	06/21/06	EPA 300.0
Nitrate (as N)	12	1	0.028	10		mg/L	N/A	06/21/06	EPA 300.0
Sulfate	73	10	0.069	10		mg/L	N/A	06/21/06	EPA 300.0
Perchlorate (1)	0.76	2.0	0.43	1	J	ug/L	N/A	06/23/06	EPA 314.0
Sulfide, Total (1)	ND	0.050	0.042	1		mg/L	06/26/06	06/26/06	EPA 376.2
Dissolved Oxygen	7.54	0.01	0.0100	1		mg/L	N/A	06/20/06	SM 4500-O G

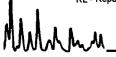
MW-7 Comment(s): (1) Results we	re evaluated to the I	Pro-Consultation of the Consultation	06-06-124 entrations >	- NECESTRAL PROPERTY OF THE	06/20/06 • MDL but •	100 PUR 100 TOTAL	100 C S SEEDS 40. (*** 100)	h a "J" flag.	
Parameter	Result	RL	MDL	DF	Qual	Units	Date Prepared	Date Analyzed	Method
Chromium, Hexavalent	1.4	0.2	0.0050	1	В	ug/L	N/A	06/20/06	EPA 218.6
Chloride	38	10	0.055	10		mg/L	N/A	06/21/06	EPA 300.0
Nitrite (as N) (1)	ND	0.10	0.015	1		mg/L	N/A	06/21/06	EPA 300.0
Nitrate (as N)	10	1	0.028	10		mg/L	N/A	06/21/06	EPA 300.0
Sulfate	67	10	0.069	10		mg/L	N/A	06/21/06	EPA 300.0
Perchlorate (1)	0.59	2.0	0.43	1	J	ug/L	N/A	06/23/06	EPA 314.0
Sulfide, Total (1)	ND	0.050	0.042	1		mg/L	06/26/06	06/26/06	EPA 376.2
Dissolved Oxygen	7.71	0.01	0.0100	1		mg/L	N/A	06/20/06	SM 4500-O G

Comment(s): (1) Results we	re evaluated to the I	MDL, conce	ntrations >	= to the	MDL but <	RL, if found	d, are qualified with	n a "J" flag.	
<u>arameter</u>	<u>Result</u>	<u>RL</u>	<u>MDL</u>	<u>DF</u>	<u>Qual</u>	<u>Units</u>	Date Prepared	Date Analyzed	<u>Method</u>
hromium, Hexavalent	1.6	0.2	0.0050	1	В	ug/L	N/A	06/20/06	EPA 218.6
hloride	41	10	0.055	10		mg/L	N/A	06/21/06	EPA 300.0
trite (as N) (1)	ND	0.10	0.015	1		mg/L	N/A	06/21/06	EPA 300.0
trate (as N)	13	1	0.028	10		mg/L	N/A	06/21/06	EPA 300.0
ulfate	71	10	0.069	10		mg/L	N/A	06/21/06	EPA 300.0
erchlorate (1)	0.75	2.0	0.43	1	J	ug/L	N/A	06/23/06	EPA 314.0
ulfide, Total (1)	ND	0.050	0.042	1		mg/L	06/26/06	06/26/06	EPA 376.2
ssolved Oxygen	7.41	0.01	0.0100	1		mg/L	N/A	06/20/06	SM 4500-O G

RL - Reporting Limit

DF - Dilution Factor ,

Qual - Qualifiers





Analytical Report



Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Date Received:

06/20/06

Work Order No:

06-06-1241

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Page 2 of 2

Client Sample Number		L	ab Sample	Numbe	r Date Collected	Matr	ix		
Method Blank					NA	" Aque	ous		
Comment(s): (1) Results were en	valuated to the N	MDL, conce	ntrations >	= to the	MDL but < f	RL, if found	d, are qualified with	h a "J" flag.	
<u>Parameter</u>	<u>Result</u>	<u>RL</u>	<u>MDL</u>	<u>DF</u>	<u>Qual</u>	<u>Units</u>	Date Prepared	Date Analyzed	<u>Method</u>
Chromium, Hexavalent (1)	0.026	0.20	0.0050	1	J	ug/L	N/A	06/20/06	EPA 218.6
Chloride (1)	ND	1.0	0.055	1		mg/L	N/A	06/20/06	EPA 300.0
Nitrite (as N) (1)	ND	0.10	0.015	1		mg/L	N/A	06/20/06	EPA 300.0
Nitrate (as N) (1)	ND	0.10	0.028	1		mg/L	N/A	06/20/06	EPA 300.0
Sulfate (1)	ND	1.0	0.069	1		mg/L	N/A	06/20/06	EPA 300.0
Perchlorate (1)	ND	2.0	0.43	1		ug/L	N/A	06/23/06	EPA 314.0
Sulfide, Total (1)	ND	0.050	0.042	1		mg/L	06/26/06	06/26/06	EPA 376.2





Quality Control - Spike/Spike Duplicate



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Date Received: Work Order No: Preparation: Method:

06/20/06 06-06-1241 EPA 3005A Filt. **EPA 6010B**

Quality Control Sample ID	Matrix	Instrument	Date Prepared		Date alyzed	MS/MSD Batch Number
MW-8	□ > Aqueou	s ICP 3300	06/21/06	06	/22/06	- 060621S06
<u>Parameter</u>	MS %REC	MSD %REC	%REC CL	<u>RPD</u>	RPD CI	Qualifiers
Antimony	106	108	72-132	2	0-10	
Arsenic	106	108	80-140	2	0-11	
Barium	104	105	87-123	1	0-6	
Beryllium	103	105	89-119	2	0-8	
Cadmium	102	105	82-124	2	0-7	
Chromium	103	105	86-122	2	0-8	
Cobalt	104	106	83-125	2	0-7	
Copper	80	82	78-126	3	0-7	
Lead	101	103	84-120	2	0-7	
Molybdenum	103	105	78-126	2	0-7	
Nickel	100	102	84-120	2	0-7	
Selenium	104	107	79-127	2	0-9	
Silver	104	106	86-128	2	0-7	
Thallium	89	91	79-121	3	0-8	
Vanadium	104	105	88-118	2	0-7	
Zinc	96	98	89-131	2	0-8	



nvironmental

Quality Control - Spike/Spike Duplicate



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

aboratories, Inc.

Date Received: Work Order No: Preparation: Method: 06/20/06 06-06-1241 EPA 3005A Filt. EPA 6010B

Quality Control Sample ID	Matrix	Instrument	Date Prepared]	Date Analyzed	MS/MSD Batch Number	
MW-8	Aqueou	s : ICP 3300	06/21/06		06/22/06	060621506	
<u>Parameter</u>	MS %REC	MSD %REC	%REC CL	RPD	RPD CL	Qualifiers	
Calcium	4X	4X	77-113	4X	0-11	Q	
Magnesium	4X	4X	56-140	4X	0-11	Q	
Potassium	105	101	83-131	2	0-7		
Sodium	4X	4X	73-127	4X	0-9	Q	

Quality Control - Spike/Spike Duplicate



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Date Received: Work Order No: Preparation: Method:

06/20/06 06-06-1241 EPA 3005A Filt. **EPA 200.8**

Quality Control Sample ID	Matrix	Matrix Instrument		1 A	Date I nalyzed	MS/MSD Batch Number
06-06-1183-4	Aquec	pus⊯∉⊸ ICP/MS A	06/21/06		06/21/06	060621S01
<u>Parameter</u>	MS %REC	MSD %REC	%REC CL	<u>RPD</u>	RPD CL	Qualifiers
Iron Manganese	142 100	134 100	80-120 80-120	6 0	0-20 0-20	3

Quality Control - Spike/Spike Duplicate



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: 06/20/06 06-06-1241 EPA 7470A Total EPA 7470A

Project BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID	Matrix	Instrument	Date Prepare	d /	Date Analyzed	MS/MSD Batch Number
06-06-1258-1	Aqueou	s Mercury	06/21/06		06/22/06	060621504
<u>Parameter</u>	MS %REC	MSD %REC	%REC CL	RPD	RPD CL	Qualifiers
Mercury	108	112	71-134	3	0-14	

RPD - Relati

Quality Control - Spike/Spike Duplicate



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: 06/20/06 06-06-1241 EPA 3520B EPA 8270C(M) Isotope Dilution

Project BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID	Matrix	Instrument	Date Prepared		Date nalyzed	MS/MSD Batch Number
06-06-1183-4	Aqueou	ıs GC/MS J	06/22/06	. 00	5/28/06	060622S01D
<u>Parameter</u>	MS %REC	MSD %REC	%REC CL	RPD	RPD CL	Qualifiers
1,4-Dioxane	83	85	50-130	3	0-20	

RPD - Relai

Quality Control - Spike/Spike Duplicate



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300

Pasadena, CA 91107-6024

Date Received: Work Order No: Preparation: Method: 06/20/06 06-06-1241 EPA 5030B EPA 8260B

Project BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID	Matrix	Instrument	Date Prepared	Α	Date nalyzed	MS/MSD Batch Number
06-07-0054-4	Aqueou	GC/MS/Z		0	7/03/06	060703501
<u>Parameter</u>	MS %REC	MSD %REC	%REC CL	<u>RPD</u>	RPD CL	Qualifiers
Benzene	96	96	88-118	0	0-7	
Carbon Tetrachloride	113	115	67-145	2	0-11	
Chlorobenzene	98	97	88-118	1	0-7	•
1,2-Dichlorobenzene	97	96	86-116	1	8-0	
1,1-Dichloroethene	. 90	93	70-130	4	0-25	
Toluene	98	98	87-123	0	0-8	
Trichloroethene	95	96	79-127	1	0-10	
Vinyl Chloride	83	85	69-129	2	0-13	
Methyl-t-Butyl Ether (MTBE)	91	94	71-131	2	0-13	
Tert-Butyl Alcohol (TBA)	7 7	. 81	36-168	6	0-45	
Diisopropyl Ether (DIPE)	101	104	81-123	3	0-9	
Ethyl-t-Butyl Ether (ETBE)	89	92	72-126	3	0-12	•
Tert-Amyl-Methyl Ether (TAME)	98	101	72-126	3 -	0-12	
Ethanol	69	70	53-149	2	0-31	

RPD - Relative Percent Difference 7440 Lincoln

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Quality Control - Spike/Spike Duplicate



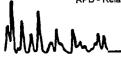
aboratories, Inc.

Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Date Received: Work Order No: Preparation: Method:

06/20/06 06-06-1241 **EPA 5030B** SRL 524M-TCP

Quality Control Sample ID	Matrix	Instrument	Date Prepare		Date I alyzed	MS/MSD Batch Number
06-06-1183-4	*Aqueou	s GC/MS M	06/21/06	₩ [*] 06	J22/06 "	060621S02
<u>Parameter</u>	MS %REC	MSD %REC	%REC CL	RPD	RPD CL	Qualifiers
1,2,3-Trichloropropane 1,4-Dioxane	66 90	66 89	80-120 80-120	0 2	0-20 0-20	3



Quality Control - Spike/Spike Duplicate



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300

Date Received: Work Order No:

N/A 06-06-1241

Pasadena, CA 91107-6024

Matrix: Aqueous										\$ 1
<u>Parameter</u>	<u>Method</u>	Quality Control Sample ID	<u>Date</u> <u>Analyzed</u>	<u>Date</u> Extracted	MS% REC	MSD % REC	%REC CL	RPD	RPD CL	Qualifiers
Chloride	EPA 300.0	06-06-1183-4	06/20/06	N/A	95	96	56-134	1	0-3	
Nitrite (as N)	EPA 300.0	06-06-1183-4	06/20/06	N/A	99	95	68-122	4	0-8	
Nitrate (as N)	EPA 300.0	06-06-1183-4	06/20/06	N/A	93	96	58-142	4	0-6	
Sulfate	EPA 300.0	06-06-1183-4	06/20/06	N/A	104	102	49-133	2	0-3	
Chromium, Hexavalent	EPA 218.6	06-06-1227-1	06/20/06	N/A	101	103	85-121	1	0-4	
Perchlorate	EPA 314.0	06-06-1183-4	06/23/06	N/A	97	97	80-120	0	0-15	



Quality Control - Duplicate



Tetra Tech, Inc.

Date Received:

N/A

3475 East Foothill Blvd., Suite 300

Work Order No:

06-06-1241

Pasadena, CA 91107-6024

Matrix: Aqueous	S. 1975 S. 20							
<u>Parameter</u>	<u>Method</u>	QC Sample ID	<u>Date Analyzed</u>	Sample Conc	DUP Conc	RPD	RPD CL	Qualifiers
Dissolved Oxygen Sulfide, Total	SM 4500-O G EPA 376.2	MW-3 06-06-1258-3	06/20/06 06/26/06	7.41 ND	7.47 ND	1 NA	0-25 0-25	

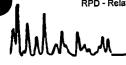
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nvironmental Quality Control - Laboratory Control Sample aboratories, Inc.



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: N/A 06-06-1241 EPA 3005A Filt. EPA 6010B

Quality Control Sample ID	Matrix	Instrument [Lab File ID	LCS Batch Number	
097-01-003-6,222	Aqueous	ICP 3300	06/22/06	060621-I-06	2 060621	L06F
Parameter	Conc	Added	Conc Recovered	LCS %Rec	%Rec CL	Qualifiers
Antimony	0.	500	0.490	98	80-120	
Arsenic	0.9	500	0.488	98	80-120	
Barium	0.9	500	0.507	101	80-120	
Beryllium	0.	500	0.487	97	80-120	
Cadmium	0.:	500	0.514	103	80-120	
Chromium	0.:	500	0.506	101	80-120	
Cobalt	0.	500	0.525	105	80-120	
Copper	0.	500	0.459	92	80-120	
Lead	. 0.	500	0.506	101	80-120	
Molybdenum	0.:	500	0.493	99	80-120	
Nickel	0.	500	0.523	105	80-120	
Selenium	0.	500	0.463	93	80-120	
Silver	0.:	250	0.242	97	80-120	
Thallium	0.8	500	0.457	91	80-120	
Vanadium	0.5	500	0.492	98	80-120	
Zinc	0.:	500	0.533	107	80-120	



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Tetra Tech, Inc.

3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

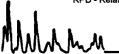
Date Received: Work Order No: Preparation:

N/A 06-06-1241 EPA 3005A Filt.

Method:

EPA 6010B

Quality Control Sample ID	Matrix	Instrument	Date Analyzed	Lab File	ID LCS E	atch Number
097-01-003-6,222	Aqueous	ICP 3300	06/22/06	060621-1	-06 ,	0621L06F
<u>Parameter</u>		Conc Added	Conc Recovered	LCS %Rec	%Rec CL	Qualifiers
Calcium	•	0.500	0.507	101	80-120	
Magnesium		0.500	0.488	98	80-120	
Potassium		5.00	5.06	101	80-120	
Sodium		5.00	5.17	103	80-120	



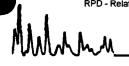


Quality Control - LCS/LCS Duplicate



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: N/A 06-06-1241 EPA 3005A Filt. EPA 200.8

Quality Control Sample ID	Carlos Francis (note 100) and Court Court	Instrument	Date Prepared	Date Analyzed	LCS/LCSD Batch Number	, \$28855283
099-10-008-738	Aqueous I	CP/MS A	06/21/06	06/21/06	060621L01F	
Parameter	LCS %REC	C LCSD %	REC %REC	CL RPD	RPD CL C	ualifiers
Iron	102	102	85-1	15 0	0-20	
Manganese	98	98	85-1	15 0	0-20	





Quality Control - LCS/LCS Duplicate



and real and real control of the

Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: N/A 06-06-1241 EPA 7470A Filt. EPA 7470A

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID	Matrix In	nstrument	Date Prepared	Date Analyzed	LCS/LCSD Bate Number	
099-04-008-2,530	Aqueous l	Mercury	06/21/06	06/21/06	060621L04F	
<u>Parameter</u>	LCS %REC	LCSD %F	REC %REC	CL RPD	RPD CL	Qualifiers
Mercury	97	97	90-12	22 0	0-14	

Mulum_

N/A



Quality Control - LCS/LCS Duplicate



aboratories, Inc.

Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Date Received: Work Order No: Preparation:

06-06-1241 **EPA 3520B**

Method:

EPA 8270C(M) Isotope Dilution

Quality Control Sample ID	Matrix In	strument	Date Prepared	Date Analyzed	LCS/LCSD Bate Number	h
099-09-004-597 A	queous G	C/MS J	06/22/06	06/27/06	060622L01D	
<u>Parameter</u>	LCS %REC	LCSD %R	EC %REC C	CL RPD	RPD CL	Qualifiers
1,4-Dioxane	83	88	50-130	5	0-20	





Quality Control - LCS/LCS Duplicate



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: N/A 06-06-1241 EPA 3520B EPA 1625CM

Project: BOU Groundwater Monitoring 2006 (PAC Wells) / 17653-0602

Quality Control Sample ID	Matrix Instrument		Date Prepared	Date Analyzed	LCS/LCSD Bato Number	zh
099-07-027-251 A	Aqueous	GC/MS H	06/23/06	06/27/06	* • 060623L09	6.90
<u>Parameter</u>	LCS %RE	C LCSD %	REC %REC	CL RPD	RPD CL	Qualifiers
N-Nitrosodimethylamine	78	80	50-1	30 4	0-20	

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alscience nvironmental Quality Control - Laboratory Control Sample aboratories, Inc.



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024 Date Received: Work Order No: Preparation: Method: N/A 06-06-1241 EPA 5030B EPA 8260B

Quality Control Sample ID	Matrix	Instrument	Date Analyze	ed Lab File ID	LCS	LCS Batch Number		
099-10-006-18,441	Aqueous	GC/MS Z	07/03/0	6 03JUL005.n		060703L01		
<u>Parameter</u>	Conc	Added	Conc Recovered	LCS %Rec	%Rec CL	Qualifiers		
Benzene	5	0	47	94	84-120			
Carbon Tetrachloride	5	0	57	114	63-147			
Chlorobenzene	5	0	48	95	89-119			
1,2-Dichlorobenzene	5	0	48	96	89-119			
1,1-Dichloroethene	5	0	46	93	77-125			
Toluene	5	0	48	95	83-125			
Trichloroethene	5	0	48	97	89-119			
Vinyl Chloride	5	0	43	87	63-135			
Methyl-t-Butyl Ether (MTBE)	5	0	47	94	82-118			
Tert-Butyl Alcohol (TBA)	2	50	210	84	46-154			
Diisopropyl Ether (DIPE)	5	0	53	107	81-123			
Ethyl-t-Butyl Ether (ETBE)		0	49	99	74-122			
Tert-Amyl-Methyl Ether (TAME)		0	52	104	76-124	•		
Ethanol	5	00	350	70	60-138			





Quality Control - LCS/LCS Duplicate



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Date Received: Work Order No: Preparation: Method: N/A 06-06-1241 EPA 5030B SRL 524M-TCP

Quality Control Sample ID	_MatrixIn	strument	Date Prepared	Date Analyzed	LCS/LCSD Batc Number	h
099-10-022-240	Aqueous G	C/MS M	06/21/06	06/21/06	060621L02	
Parameter	LCS %REC	LCSD %F	EC %REC	CL RPD	RPD CL	Qualifiers
1,2,3-Trichloropropane	86	91	80-12	0 6	0-20	
1,4-Dioxane	83	87	80-12	0 4	0-20	



Quality Control - LCS/LCS Duplicate



Tetra Tech, Inc. 3475 East Foothill Blvd., Suite 300 Pasadena, CA 91107-6024

Date Received: Work Order No: N/A

06-06-1241

Matrix: Aqueous										
<u>Parameter</u>	<u>Method</u>	Quality Control Sample ID	<u>Date</u> Extracted	<u>Date</u> Analyzed	LCS % REC	LCSD % REC	%REC CL	RPD	RPD CL	<u>Qual</u>
Chloride	EPA 300.0	099-05-118-3,422	N/A	06/20/06	94	95	81-111	1	0-5	
Nitrite (as N)	EPA 300.0	099-05-118-3,422	N/A	06/20/06	92	94	73-115	1	0-26	
Nitrate (as N)	EPA 300.0	099-05-118-3,422	N/A	06/20/06	97	97	87-111	0	0-12	
Sulfate	EPA 300.0	099-05-118-3,422	N/A	06/20/06	99	97	89-107	2	0-13	
Chromium, Hexavalent	EPA 218.6	099-05-124-487	N/A	06/20/06	98	98	95-107	0	0-20	
Perchlorate	EPA 314.0	099-05-203-423	N/A	06/23/06	107	104	85-115	3	0-15	



Glossary of Terms and Qualifiers



Work Order Number: 06-06-1241

Qualifier	<u>Definition</u>
*	See applicable analysis comment.
:1	Surrogate compound recovery was out of control due to a required sample dilution, therefore, the sample data was reported without further clarification.
2	Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification.
3	Recovery of the Matrix Spike or Matrix Spike Duplicate compound was out of control due to matrix interference. The associated LCS and/or LCSD was in control and, therefore, the sample data was reported without further clarification.
4	The MS/MSD RPD was out of control due to matrix interference. The LCS/LCSD RPD was in control and, therefore, the sample data was reported without further clarification.
5	The PDS/PDSD associated with this batch of samples was out of control due to a matrix interference effect. The associated batch LCS/LCSD was in control and, hence, the associated sample data was reported with no further corrective action required.
Α	Result is the average of all dilutions, as defined by the method.
В	Analyte was present in the associated method blank.
С	Analyte presence was not confirmed on primary column.
E	Concentration exceeds the calibration range.
H	Sample received and/or analyzed past the recommended holding time.
J .	Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated.
N	Nontarget Analyte.
ND	Parameter not detected at the indicated reporting limit.
Q	Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater.
Ų	Undetected at the laboratory method detection limit.
Χ	% Recovery and/or RPD out-of-range.
Ζ .	Analyte presence was not confirmed by second column or GC/MS analysis.

	TETRA TECH, INC.
	3475 E. FOOTHILL BLVD. PASADENA, CALIFORNIA 91107 TELEPHONE (626) 351-4664
	PASADENA, CALIFORNIA 91107
	TELEPHONE (626) 351-4664
• •	CAV (COR)254 5004

SHIPPED TO: CAL SCIENC

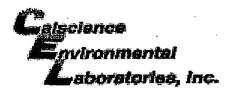
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WORK ORDER #: 06 - 0 6 - 1 2 4 1

Cooler ____ of ___

SAIVIPLE REC	CIFI FURIVI
CLIENT: Tetractech	DATE: 6/20/6
TEMPERATURE - SAMPLES RECEIVED BY:	
CALSCIENCE COURIER: Chilled, cooler with temperature blank provided. Chilled, cooler without temperature blank. Chilled and placed in cooler with wet ice. Ambient and placed in cooler with wet ice. Ambient temperature.	LABORATORY (Other than Calscience Courier): °C Temperature blank. °C IR thermometer. Ambient temperature.
°C Temperature blank.	Initial:
CUSTODY SEAL INTACT: Sample(s): Cooler: No (Not Intact)	:Not Applicable (N/A): Initial:
Chain-Of-Custody document(s) received with samples	
COMMENTS:	

APPENDIX C QA/QC SUMMARY

APPENDIX C

QUALITY ASSURANCE/QUALITY CONTROL SUMMARY

1.1 QUALITY ASSURANCE/QUALITY CONTROL SUMMARY

The Quality Assurance/Quality Control (QA/QC) Summary is the relevant QA/QC information associated with the Burbank Operational Unit sampling data set (PACWELLS). The QA/QC Summary contains the following three subjects, which are addressed in detail:

- Data validation concepts, rationale, and practices;
- Data quality objectives, evaluation, and implications; and

1.1.1 SELECTED DEFINITIONS/CRITERIA OF TERMS

1.1.1.1 Holding Times

The U. S. Environmental Protection Agency (U.S. EPA) has established maximum time intervals (holding times) between the collection, extraction, and analysis of samples. All compliant results must be obtained within holding times or the results are considered deficient. Samples analyzed outside of holding times must be qualified.

1.1.1.2 Laboratory and Field Blanks

Laboratory and field blanks are samples used to determine if environmental sample results may be positively biased by laboratory or field contamination. Laboratory blank results indicate contamination due to laboratory operations only, while field blank results indicate contamination from field and/or laboratory operations. Laboratory blanks contaminated above the Practical Quantitation Limit (PQL) indicate a need for corrective action.

1.1.1.3 Matrix Spike/Matrix Spike Duplicate (MS/MSD)

Matrix spike samples are environmental samples that are spiked with known concentrations of target analytes. The recovery of the target analytes is used to evaluate the effects of the sample matrix. Matrix effects are considered site specific. One MS/MSD sample is analyzed for every 20 environmental samples. The matrix spike duplicate results may be compared to the matrix spike results in order to determine precision.

1.1.1.4 Laboratory Control Sample (LCS)

The LCS determines if the analytical system is in control and consists of reagent grade (analyte free) water spiked with known concentrations of target analytes. Results from the LCS are considered free of any matrix effects and analyte recoveries outside control limits are used to qualify data.

1.1.1.5 Surrogates

For most methods, surrogate compounds are added to every sample at the beginning of sample preparation and are used to monitor the analytical process and give information concerning matrix effects. Surrogate recoveries are the single most useful QC entity for evaluating environmental analytical data. The ubiquitous use of surrogates in the analytical methods has afforded a large database of results from which useful correlated information can be extracted. Surrogates are chemically similar to target analytes and their

recovery within control limits indicates the process is in control. Surrogates are the primary indicators of matrix effects.

1.1.1.6 Second Column Confirmation

All organic analysis that results in analyte detection should be confirmed in order to have confidence in the result. In the case of gas chromatography/mass spectrometry (GC/MS) analysis, analyte peaks at the correct retention time are confirmed by the mass spectra. For GC or high performance liquid chromatography (HPLC) analysis, a second analytical column and/or a second detector is used for to confirm the presence of the analyte. Unless an analyte is confirmed, its presence cannot be proved.

1.1.1.7 Temperature Blanks

Temperature blanks are placed in coolers with environmental samples in order to determine the temperature of the samples when they arrive at the lab. Temperature blanks typically consist of water in a container similar to the sample containers. Upon receipt at the lab, the temperature blanks are opened and a thermometer is inserted directly into the liquid. Alternatively, the temperature of the samples is measured using an infrared thermometer. The criterion is 4 degrees Celsius, plus or minus 2 degrees. Samples that arrive at the laboratory shortly after sample collection (less than 4 hours) may not have sufficient time for temperature equilibration. In these cases, samples may exceed the upper temperature limit of 6 degrees Celsius, but must be below ambient temperatures.

1.1.1.8 Field Audits

Field audits determine if the sampling procedures used by the field crew are in accordance with standard operating procedures. The techniques used to collect the samples are analyzed to determine if the samples are being collected correctly.

1.1.1.9 Sample Delivery Group (SDG)

The SDG is a laboratory-defined collection of sample results together with the corresponding quality control results. These results are organized under a unique group heading. The laboratory determines the method of grouping the sample results under an SDG and each SDG may contain samples collected at various times and with different matrix types. Generally, each SDG consists of the results for a group of samples received by the laboratory on a single day.

1.1.1.10 Data Gaps

Data gaps may be generated by both field sampling activities and laboratory data problems. Field activities that may produce data gaps include difficulty accessing the sampling location, which results in no sample being collected, or damage and subsequent loss of samples before they reach the laboratory. Laboratory QC errors resulting in data that must be qualified as rejected will also leave data gaps in the analytical results. If necessary, data gaps may be closed quickly by resampling and reanalysis. If the results are not time critical, the gap may be closed during the next quarter of sampling.

1.1.1.11 Corrective Actions

Corrective actions are performed in response to data or conditions that are not in analytical control. Corrective actions are performed in an attempt to bring the error condition back under control. Corrective actions are documented by a corrective action report (CAR) and are included in the laboratory's SDG data package.

1.1.2 DATA VALIDATION RATIONALE AND GUIDELINES

1.1.2.1 Controlling Documents

The following documents were used for data validation.

- USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review (Publication OSWER 9240.1-05A-P, EPA-540/R-99/008, October 1999); and
- USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (Publication OSWER 9240.1-35, EPA-540/R-01/008, July 2002).

These two documents are hereinafter collectively referred to as the National Functional Guidelines.

1.1.2.2 Data Validation Theory and Matrix Effects

The practice of data validation in the environmental organic chemistry field has been the subject of debate for many years. Determining the validity of environmental data results when matrix effects are suspected is not an exact science, and professional judgment concerning matrix effects is used to help guide the data to its best logical interpretation and evaluation.

The overall QC of environmental sample analysis can be divided into two main categories. These categories are generally considered to be "method QC" and "instrument QC." Both types of QC operate independently to validate the data and qualify the results.

Instrument QC parameters are often well defined and well understood and are based on the tangible physical laws of analytical instrumentation. Instrument QC parameters have to do with (but are not limited to) the calibration, chromatography, and detection aspects of environmental data analysis. Instrument QC parameters are considered independent from a sample's matrix and/or matrix effects.

Conversely, method QC parameters do not yield results that are as well defined, since they are based in part on problems associated with the intangible and/or unknown effects of the sample matrix. Method QC parameters have to do with (but are not limited to) the spiking, extraction, and spike recovery aspects of environmental data analysis. Method QC parameters are considered dependent on a sample's matrix and/or matrix effects.

When evaluating environmental data results with pronounced or unknown matrix effects, a conservative approach to the validation is required. The method QC parameters are rigidly applied and validations are conferred to entire data sets based on one sample's bias.

1.1.2.3 Data Validation Rationale

The National Functional Guidelines were written for use with the Contract Laboratory Program (CLP) methods as outlined in the CLP Statement of Work (SOW). The SOW contains methods for volatile and semivolatile GC/MS analysis, two-column GC pesticide analysis, and inductively coupled plasma (ICP) metal analysis. These methods do not differ significantly in the application of the basic quality control parameters from those found in the corresponding SW846 methods for volatile, semivolatile, pesticide, and ICP metals analyses (hereafter referred to as the SW methods). The target compounds in the CLP are a subset of the SW846 target compounds.

Since the CLP methods and the SW846 methods have similar QC instructions, the *National Functional Guidelines* are usable for the SW methods. In order to validate analytical methods that have no corresponding CLP method validation guidance, logical extrapolations are determined by modeling the pertinent CLP rationale. The resulting validated data have a professional judgment component that allows the validation to be tailored to the individual project. Since the validation of environmental results is not an exact science, interpretive judgments are sometimes required for complex data

1.1.2.4 Validation Qualifiers

- B: The sample result is less than 5 times (10 times for common organic laboratory contaminants) the blank contamination. The result is considered not to have originated from the environmental sample, because cross-contamination is suspected.
- J: The analyte was positively identified and the result is usable; however, the analyte concentration is an estimated value.
- R: The sample result is rejected and not usable for any purpose. The presence or absence of the analyte cannot be verified.
- U: The analyte was not detected at or above the reporting detection limit (RDL).
- UJ: The analyte was not detected above the MDL; however, the MDL is uncertain and may be elevated above normal levels.
- Y: Confirmation column results indicate a non-detect for the target analyte.

1.1.2.5 Qualifier Descriptors

- a: The analyte was found in the method blank.
- b: The surrogate spike recovery was outside quality control criteria.
- c: The MS and/or MSD recoveries were outside control limits.
- d: The laboratory control sample recovery was outside control limits.
- e: A holding time violation occurred.
- f: The duplicate/replicate sample's relative percent difference (RPD) was outside the control limit.
- g: The data met prescribed criteria as detailed in the QAPP.
- h: The required second column confirmation was not performed.
- k: The analyte was found in a field blank.
- 1: The second column confirmation result indicates the analyte was not confirmed.
- n: The laboratory case narrative indicated a QC problem.
- p: Professional judgment determined the data should be qualified.
- q: The analyte detection was below the PQL.
- r: The result is above the instrument's calibration range.
- t: The temperature was outside acceptance criteria.

1.1.2.6 Level One Validation Guidelines

Organic Validation Guidelines

Sample Preservation

- As a rule, all samples are required to be preserved at a temperature of 4 degrees Celsius, plus or minus 2 degrees. Additional preservation criteria are method specific. The temperature criterion applies to all samples.
- Samples placed in a cooler and transported directly to the laboratory with short transit times (less than 4 hours) do not allow for temperature equilibration. The temperature of samples with short transit time must be below ambient temperature with evidence of cooling in progress (ice or ice-substitute present).

• Samples with temperatures in excess of six degrees Celsius but less than or equal to 12 degrees Celsius are qualified J for detected analytes and UJ for non-detects.

- Samples in gross excess (>12 degrees) of the temperature criteria are qualified **J** for detected analytes and non-detects are qualified **R**.
- The descriptor t is used to indicate sample temperature qualification.

Holding Times

- For volatile organic analyses (VOA) samples, analysis after 14 days (7 days if not pH preserved) from collection are qualified **J** and **UJ**.
- For semivolatile (SV) samples, water samples extracted after 7 days (14 days for soil) are
 qualified J and UJ. Samples analyzed after 40 days from extraction are also qualified J
 and UJ.
- If holding times are grossly exceeded (greater than 2 times the normal holding time), then positive results are qualified **J** and non-detects are qualified **R**.
- The descriptor **e** is used to denote holding time violations.

Blanks

- Analytes found in associated environmental samples at or below 5 times (10 times common organic analytes) of the method or field blank analyte concentrations are qualified **B**.
- The descriptor **a** is used to indicate method blank contamination.
- The descriptor k is used to indicate field blank contamination.

Surrogates

- For VOA (GC/MS) samples, there are three cases. Any single surrogate failure will cause qualification.
 - Case #1: Recovery above upper limit, then **J** qualify detected analytes. Do not qualify non-detected analytes.
 - Case #2: Recovery between lower limit and 10 percent, then J and UJ.
 - Case #3: Recovery below 10 percent, then J positive results and R non-detects.
- For SV (GC/MS) samples, there are four cases. Except for case four, two surrogate failures (within each fraction) will cause fraction specific qualification.
 - Case #1: Recovery above upper limit, then J only. No UJ.

- Case #2: Recovery between lower limit and 10 percent, then J and UJ.
- Case #3: Recovery of one surrogate above upper limit and one surrogate below the lower limit but above 10 percent, then qualify as in case #2.
- Case #4: Any one surrogate below 10 percent, then J positive results and R non-detects.
- For SV (GC) samples.
 - Case #1: Recovery above upper limit, then **J** only positive results. Non-detects are not qualified.
 - Case #2: Recovery between lower limit and 10 percent, then J positive results. Non-detects are qualified UJ.
 - Case #3: Recovery below 10 percent, then J positive results and R non-detects.
- The descriptor **b** is used to indicate surrogate failure qualification.

Laboratory Control Sample

- For laboratory control sample (LCS) qualifications, the specific analytes spiked into the LCS sample must always be qualified. All target analytes are spiked into the LCS.
- For all methods requiring LCS recoveries there are 2 cases.
 - Case #1: LCS recovery is above upper limit, then **J** detected analytes only. Do not qualify non-detects.
 - Case #2: LCS recovery is below lower limit then J positive results and R non-detects.
- The descriptor **d** is used to indicate LCS qualification.

Matrix Spike/Matrix Spike Duplicates

- The target analytes spiked into the MS/MSD are listed in the project specific QAPP.
- There are two cases for qualification based on the MS/MSD results.
 - Case #1: Non-compliant spike recoveries comprise the first case for qualification based on MS/MSD results. MS and MSD spike recoveries outside of control limits, where the LCS demonstrates that the analytical system was in control, are attributed to the effects of the sample matrix. If both the MS and MSD fail spike recovery criteria as indicated below, qualify based on the least compliant recovery.

- Recovery above upper limit, then **J** detected compounds only. Do not qualify non-detects.

- Recovery between lower limit and 10 percent, then J detected compounds and UJ non-detects.
- Recovery below 10 percent, then J detected compounds and R non-detects.
- Case #2: Non-compliance of the RPD value is the second case for qualification of data based on the MS/MSD results. MS/MSD RPDs are calculated from the analyte concentrations of the MS and MSD. If the RPD is outside the control limit, the precision is in question, and the accuracy is compromised.
 - RPD outside the control limit, then qualify the related samples with J for detected compounds and UJ non-detects.
- The descriptor c is used to indicate MS/MSD qualification based on the percent recovery of the spiked analytes.
- The descriptor **f** is used to indicate RPD failure.

Second Column Confirmation

For certain GC or HPLC methods, second column/detector confirmation is required for detected analytes. Refer to the relevant QAPP for method and analyte specific requirements.

Second column results are used to <u>confirm the actual presence or absence of a target analyte</u>. U.S. EPA guidelines state "If the qualitative criteria for both columns were <u>not</u> met, all target compounds that are reported detected should be considered non-detected." Therefore, any compound detection on only one column is not considered a target compound hit.

- For the situation where a compound was detected on the primary column and not detected on the confirmation column, consider the value reported to be not detected. Qualify the result with Y and use the I descriptor.
- In the case of a detection on the primary column where the required second column confirmation was not performed, then qualify the result with **R** and use the **h** descriptor.

Field Duplicate Samples

Field duplicate samples are collected to assess the precision of the sample collection and laboratory analytical process. As a rule, both the sample and its duplicate result must be at or above the PQL in order to calculate a meaningful RPD and if both results are below the PQL the RPD is not calculated. However, if one result is below the PQL (assume zero for a non-detect) and the other result

significantly above (10 times) the PQL a RPD is calculated. If the RPD is outside the control limit, the precision is in question, and the accuracy is compromised. The qualification resulting from the sample and its duplicate non-compliant RPD apply only to the sample and it's duplicate and is analyte specific.

- If the RPD is outside the control limit, then qualify the sample and its duplicate with J for detected compounds and UJ non-detects.
- The descriptor **f** is used to indicate RPD failure.

Inorganic Validation Guidelines

Sample Preservation

- As a rule, all samples are required to be preserved at a temperature of 4 degrees Celsius, plus or minus 2 degrees. Additional preservation criteria are method-specific. The temperature criterion applies to all samples except ICP metals and mercury in a water matrix, which are exempt from temperature preservation.
- Samples placed in a cooler and transported directly to the laboratory with short transit times (less than 4 hours) do not allow for temperature equilibration. The temperature of samples with short transit time must be below ambient temperature with evidence of cooling in progress (ice or ice-substitute present).
- Samples with temperatures in excess of six degrees Celsius but less than or equal to 12 degrees Celsius are qualified J for detected analytes and UJ for non-detects.
- Samples in gross excess (more than 12 degrees) of the temperature criteria are qualified J
 for detected analytes and non-detects are qualified R.
- The descriptor **t** is used to indicate sample temperature qualification.

Holding Times

- Holding times are measured from the sampling date.
- Holding times for inorganic compounds vary from 24 hours for analyses such as chromium
 VI and pH to six months for ICP metals. Results produced from analyses performed
 beyond the holding time are qualified as estimated J for detected values and UJ for
 nondetects.
- If holding times are grossly exceeded (greater than 2 times the normal holding time), then positive results are qualified **J** and non-detects are qualified **R**.
- The descriptor e is used to denote holding time violations.

Blanks

• Equipment blanks and/or laboratory blanks are evaluated for contaminants.

- Analytes found in associated environmental samples at or below 5 times the blank analyte contamination are qualified **B**.
- Analytes qualified for laboratory blank contamination are denoted with a descriptor a.
- Analytes qualified for equipment blank contamination are denoted with a descriptor k.

Laboratory Control Sample

- For LCS qualifications, the specific analytes spiked into the LCS sample must always be qualified. All target analytes are spiked into the LCS.
- LCS recovery is above upper limit then J detected analytes only. Do not qualify non-detects.
- LCS recovery is below lower limit then J positive results and R non-detects.
- Analytes qualified for LCS failure are denoted with a descriptor d.

Matrix Spike/Matrix Spike Duplicate

The target analytes spiked into the MS/MSD are listed in the project specific QAPP. Each specific MS or MSD spiking analyte that fails recovery criteria produces qualification of the matching analyte in the site associated environmental samples. Where both the MS and MSD fail criteria, qualify based on the least compliant recovery.

- MS/MSD recovery results are not used for qualification if the analyte concentration in the environmental sample used for the MS/MSD exceeds the spike concentration by a factor of 4 or more.
- If the MS and/or MSD recovery exceed the upper control limit, then J detected compounds only. Do not qualify non-detected compounds.
- If the MS and/or MSD recovery falls between the lower limit and 10 percent, then J detected compounds and UJ non-detects.
- If the MS or MSD recovery is less than 10 percent, then J detected analytes and R non-detected analytes.
- The descriptor c is used to indicate MS/MSD qualification based on the percent recovery of the spiked analytes.
- MS/MSD RPDs are calculated from the analyte concentrations of the MS and MSD.
 If the RPD is outside the control limit, the precision is in question, and the accuracy is compromised.
- MS/MSD RPD results are not used for qualification if the analyte concentration in the

environmental sample used for the MS/MSD exceeds the spike concentration by a factor of 4 or more.

- RPD outside the control limit, then qualify the related sample results with **J** for detected compounds and **UJ** non-detects.
- The descriptor f is used to indicate RPD failure.

Field Duplicate Samples

Field duplicate samples are collected to assess the precision of the sample collection and laboratory analytical process. As a rule, both the sample and its duplicate result must be at or above the PQL in order to calculate a meaningful RPD and if both results are below the PQL, the RPD is not calculated. However, if one result is below the PQL (assume zero for a non-detect) and the other result significantly above (10 times) the PQL a RPD is calculated. If the RPD is outside the control limit, the precision is in question, and the accuracy is compromised. The qualification resulting from the sample and its duplicate non-compliant RPD apply only to the sample and it's duplicate and is analyte specific.

- If the RPD is outside the control limit, then qualify the sample and its duplicate with J for detected compounds and UJ non-detects.
- The descriptor **f** is used to indicate RPD failure.

1.1.3 SUMMARY OF DATA QUALITY OBJECTIVES AND COMPLIANCE

1.1.3.1 Data Quality Objectives

Data quality objectives (DQOs) are qualitative and quantitative statements developed by data users to specify the quality of data from field and laboratory data collection activities. These DQOs must be carefully designed to support specific decisions or regulatory actions. The DQOs describe which data are needed, why the data are needed, and how the data will be used to address the problem being investigated. DQOs also establish numeric limits for the data to allow the data user to determine whether the data collected are of sufficient quality for use in their intended application.

The usability of the data collected during this investigation depends on its quality. A number of factors relate to the quality of data, and sample collection methods are as important to consider as methods used for sample analysis. Following standard operating procedures for both sample collection and analysis reduces sampling and analytical error. Complete chain-of-custody documentation and adherence to required sample preservation techniques, holding times and proper shipment methods ensure sample integrity. Obtaining valid and comparable data also requires adequate QA/QC procedures and documentation, as well as established detection and control limits.

Quantitation limits are based on the extent to which the field equipment, laboratory equipment, or analytical process can provide accurate measurements of consistent quality for specific constituents in field samples. The quantitation limit for a given analysis will vary depending on instrument sensitivity and matrix effects.

1.1.3.2 Precision, Accuracy, Completeness, and Comparability

The effectiveness of a QA program is measured by the quality of data generated by the laboratory. Data quality is judged in terms of its precision, accuracy, completeness, and comparability. These terms are described as follows:

Accuracy

Accuracy is the degree of agreement of a measurement or average of measurements with an accepted reference or "true" value, and is a measure of bias in the system. The accuracy of a measurement system is impacted by the errors introduced through the sampling process, field contamination, preservation, handling, sample matrix, sample preparation, and analytical techniques.

For this project, laboratory accuracy of the measurement data will be assessed and controlled. Results for blanks, matrix spikes, LCS, and surrogates will be the primary indicators of accuracy. These results will be used to control accuracy by requiring that they meet specified criteria. As spiked samples are analyzed, spike recoveries will be calculated and compared to pre-established acceptance limits.

Acceptance limits are based upon previously established laboratory performance for similar samples. In this approach, the control limits reflect the minimum and maximum recoveries expected for individual measurements for an in-control system. Recoveries outside the established limits indicate some assignable cause, other than normal measurement error, and possible need for corrective action. This includes recalibration of the instrument, reanalysis of the QC sample, reanalysis of the samples in the batch, or flagging the data as suspect if the problems cannot be resolved. For contaminated samples, recovery of matrix spikes may depend on sample homogeneity, matrix interference, and dilution requirements for quantification.

Precision

Precision is a measure of agreement among individual measurements of the same property under prescribed similar conditions. When control limits are established for accuracy, it automatically identifies the precision of the method. In the analysis of samples in a preparation batch, if the recoveries of analytes in the LCS are within the control limits, then the precision is also within limits.

Precision is also determined from duplicate sample analysis and MS/MSD analysis. The precision is quantified by the RPD value calculated from the duplicate results.

Completeness

Completeness is a measure of the amount of valid data obtained from a measurement system compared to the amount expected to be obtained under correct, normal conditions.

Successful analyses are defined as those where the samples arrived at the laboratory intact, properly preserved, in sufficient quantity to perform the requested analyses, and accompanied by a completed chain-of-custody. Furthermore, the sample must be analyzed within the specified holding time and in such a manner that analytical QC criteria described in this document are met.

Factors that adversely affect completeness include:

• Receipt of samples in broken containers;

 Receipt of samples in which chain of custody or sample integrity is compromised in some way;

• Samples received with insufficient volume to perform initial analyses or repeat analyses, if initial efforts do not meet QC acceptance criteria;

Improperly preserved samples; and

 Samples held in the field or laboratory longer than expected, thereby jeopardizing holding time requirements.

Completeness for the entire project also involves completeness of field and laboratory documentation, whether all samples and analyses specified in the Sampling and Analysis Plan have been processed, and whether the procedures specified in the SAP, Work Plan, and Laboratory Standard Operating Procedures (SOPs) have been implemented.

Comparability

Comparability expresses the confidence with which one data set can be compared to another data set measuring the same property. Comparability is ensured through the use of established and approved sample collection techniques and analytical methods, consistency in the basis of analysis (wet or dry weight, volume, etc.), consistency in reporting units, and analysis of standard reference materials.

1.1.3.3 Specific Measurement DQOs for Evaluating Data DQO Compliance

1. Precision is expressed in RPD values. Spiked (MS/MSD) and unspiked duplicate field samples are analyzed in order to determine precision.

2. Accuracy is expressed as a percentage of the data outside the QC entity's control limits. The percent recoveries from laboratory control sample spikes, matrix spikes and surrogate spikes are used to determine accuracy.

The samples for this data set were examined to determine compliance with the DQOs. The results are listed below.

The following methods analyzed samples for the BOU PACWELLS project and resulted in usable data of known precision and accuracy except as listed below. Several analytes had detections below the PQL and are defined as an estimated value. All of these detections are usable data.

Method 314.0 for Perchlorate

No adverse QC issues were detected.

Method 1625C (M) for low level N-Nitrosodimethylamine

No adverse QC issues were detected.

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Method 8270C (M) for low level 1,4-Dioxane

No adverse QC issues were detected

Method 524.2 for 1,2,3-Trichloropropane

Low matrix spike recovery values qualified as estimated 1,2,3-Trichloropropane in sample MW-4. The estimated data is usable for its intended purpose.

Method SW6010B/SW7470A for Title 22 Metals

Low matrix spike recovery values qualified as estimated Copper in sample MW-4. The estimated data is usable for its intended purpose.

See Section 1.1.3.7 Blank Contamination below

Method 200.8 for Iron and Manganese

Low matrix spike recovery values qualified as estimated Iron in sample MW-4. The estimated data is usable for its intended purpose.

Method 218.6 for Hexavalent Chromium

No adverse QC issues were detected.

Method SW8260B for Volatile Organic Compounds

See Section 1.1.3.7 Blank Contamination below

Method 300.0 for Common Inorganic Ions

No adverse QC issues were detected.

Method 376.2 for Sulfide

No adverse QC issues were detected.

Method SM 4500-O G for Dissolved Oxygen

No adverse QC issues were detected

1.1.3.4 Completeness

The completeness of this data set was above the DQO criterion of 90 percent. The DQO was satisfied.

1.1.3.5 Data Gaps

All data are usable for their intended purpose except for that data qualified due to method blank contamination. The data qualified for method blank contamination is less than four percent of the total data. No significant data gaps exist.

1.1.3.6 Holding Times Compliance

All samples were analyzed within method specified holding times.

1.1.3.7 Blank Contamination

The method blank the laboratory analyzes is a QC sample that determines if laboratory operations have introduced contamination into the analytical process. An analyte free method blank indicates laboratory operations are not introducing contamination into the process. Conversely, analyte detections in the method blank indicate laboratory sponsored detections. Similar detections in environmental samples that match method blank contamination are qualified with a "B" qualifier.

Detected results with a "B" qualifier indicate the results are due to laboratory sponsored activities and were not present in the native sample. Because the detection was likely caused by laboratory cross-contamination the detected result is considered not usable.

Method SW6010B samples MW-4, MW-4D, MW-5, and MW-6 were qualified due to method blank contamination for analytes Selenium, Thallium, and Zinc. The data is generally not usable.

Method SW6010B samples MW-3, MW-7, and MW-8 were qualified due to method blank contamination for analytes Chromium, Cobalt, and Nickel. The data is generally not usable

Method SW8260B samples MW-5, MW-4D, MW-3, MW-7, and MW-8 were qualified due to method blank contamination for analytes Acetone and Methylene Chloride. The data is generally not usable.

1.1.3.8 Other QC Problems

None to report.